

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG101720\
 Data File : BG046968.D
 Acq On : 18 Oct 2020 10:26
 Operator : CG/JU
 Sample : L4201-13
 Misc : GCMS CONFIRMATION
 ALS Vial : 34 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 C0AP3

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-BG101520MA.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	3.371	3	5	7	rBV2	6969	6300	0.61%	0.081%
2	3.406	7	11	21	rVB	659942	721177	70.10%	9.284%
3	3.476	21	23	26	rVB	1515	1506	0.15%	0.019%
4	3.723	61	65	67	rBV	5703	7571	0.74%	0.097%
5	3.752	67	70	74	rVB	21724	25348	2.46%	0.326%
6	4.058	120	122	125	rBV	1360	1670	0.16%	0.021%
7	4.134	131	135	137	rBV	4826	6032	0.59%	0.078%
8	4.234	147	152	156	rVB	16222	22357	2.17%	0.288%
9	4.346	167	171	176	rBV	2536	4034	0.39%	0.052%
10	4.422	179	184	187	rBV2	1726	2924	0.28%	0.038%
11	4.516	197	200	202	rBV	1857	2052	0.20%	0.026%
12	4.787	241	246	251	rBV3	3003	5137	0.50%	0.066%
13	4.992	279	281	284	rVB	2242	1986	0.19%	0.026%
14	5.086	292	297	298	rBV2	4403	5643	0.55%	0.073%
15	5.198	315	316	322	rVB2	1542	2133	0.21%	0.027%
16	5.251	322	325	328	rBV2	1725	1736	0.17%	0.022%
17	5.562	372	378	385	rVB	51394	73394	7.13%	0.945%
18	5.703	397	402	405	rVV2	2425	2953	0.29%	0.038%
19	6.062	461	463	468	rBV3	2398	3569	0.35%	0.046%
20	6.150	475	478	480	rBV3	1335	1500	0.15%	0.019%
21	6.332	506	509	512	rVB3	2006	2226	0.22%	0.029%
22	6.479	532	534	537	rVB	1420	1384	0.13%	0.018%
23	6.549	539	546	551	rVV2	7395	11669	1.13%	0.150%
24	6.743	575	579	582	rVB	1056	1562	0.15%	0.020%
25	6.808	587	590	596	rVB	1474	2473	0.24%	0.032%
26	7.043	626	630	633	rBV2	1659	2595	0.25%	0.033%
27	7.113	639	642	644	rBV2	1217	1587	0.15%	0.020%
28	7.190	654	655	658	rBV2	1590	1419	0.14%	0.018%
29	7.290	670	672	675	rBV	1303	1788	0.17%	0.023%
30	7.971	785	788	791	rVB	1636	1735	0.17%	0.022%
31	8.065	795	804	815	rBV	249258	413985	40.24%	5.329%
32	8.188	821	825	827	rBV	2010	2808	0.27%	0.036%
33	8.247	830	835	837	rVB2	1592	1382	0.13%	0.018%
34	8.435	865	867	872	rVB2	1768	2236	0.22%	0.029%

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35	8.559	885	888	890	rVB2	1364	1423	0.14%	0.018%
36	8.600	894	895	900	rBV2	918	1467	0.14%	0.019%
37	8.711	910	914	915	rBV2	1278	1558	0.15%	0.020%
38	8.823	929	933	938	rBV2	1009	1650	0.16%	0.021%
39	9.675	1074	1078	1080	rBV	1117	1561	0.15%	0.020%
40	10.262	1176	1178	1181	rVB	2181	1714	0.17%	0.022%
41	10.879	1275	1283	1291	rBV	304981	543820	52.86%	7.000%
42	10.979	1297	1300	1303	rBV	1012	1545	0.15%	0.020%
43	11.285	1349	1352	1353	rBV	2089	1487	0.14%	0.019%
44	11.943	1462	1464	1470	rVB2	1178	1528	0.15%	0.020%
45	12.055	1477	1483	1487	rVB2	1303	3396	0.33%	0.044%
46	12.489	1555	1557	1560	rBV	1669	1694	0.16%	0.022%
47	13.300	1691	1695	1698	rVB	1446	2190	0.21%	0.028%
48	13.341	1698	1702	1703	rBV2	1139	1409	0.14%	0.018%
49	13.435	1715	1718	1720	rBV	1741	1627	0.16%	0.021%
50	13.506	1727	1730	1733	rBV2	1409	2005	0.19%	0.026%
51	13.847	1783	1788	1789	rVB2	1024	1411	0.14%	0.018%
52	14.205	1847	1849	1852	rBV2	1286	1471	0.14%	0.019%
53	14.246	1855	1856	1859	rBV	1558	1781	0.17%	0.023%
54	14.305	1862	1866	1867	rBV	1254	1426	0.14%	0.018%
55	14.375	1876	1878	1882	rBV	1151	1560	0.15%	0.020%
56	14.458	1890	1892	1897	rVB	1819	1835	0.18%	0.024%
57	14.693	1924	1932	1940	rBV2	473329	760156	73.89%	9.785%
58	15.092	1996	2000	2003	rVB4	2474	3302	0.32%	0.043%
59	15.169	2011	2013	2016	rVB2	2347	2350	0.23%	0.030%
60	15.251	2025	2027	2030	rBV	2017	2216	0.22%	0.029%
61	15.351	2043	2044	2049	rBV2	1221	1397	0.14%	0.018%
62	15.415	2054	2055	2060	rVB3	1896	1849	0.18%	0.024%
63	15.474	2060	2065	2068	rBV2	2407	3063	0.30%	0.039%
64	15.539	2072	2076	2079	rVB3	3163	4653	0.45%	0.060%
65	15.621	2087	2090	2092	rVB3	1105	1425	0.14%	0.018%
66	15.733	2106	2109	2110	rBV2	3066	3164	0.31%	0.041%
67	15.744	2110	2111	2117	rVB2	2897	3285	0.32%	0.042%
68	15.797	2119	2120	2122	rBV	2196	1577	0.15%	0.020%
69	15.856	2128	2130	2132	rVB	1824	1558	0.15%	0.020%
70	15.956	2145	2147	2148	rBV	2429	1653	0.16%	0.021%
71	16.032	2158	2160	2166	rVB3	2105	3088	0.30%	0.040%

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72	16.097	2168	2171	2173	rVB3	2383	2691	0.26%	0.035%
73	16.126	2173	2176	2178	rBV	1725	1785	0.17%	0.023%
74	16.238	2190	2195	2198	rBV5	8878	11195	1.09%	0.144%
75	16.285	2200	2203	2206	rVV4	2621	3485	0.34%	0.045%
76	16.308	2206	2207	2210	rVB3	2325	1882	0.18%	0.024%
77	16.373	2216	2218	2219	rBV2	1932	1398	0.14%	0.018%
78	16.391	2219	2221	2224	rVV3	4279	5998	0.58%	0.077%
79	16.438	2224	2229	2240	rVB	57046	97915	9.52%	1.260%
80	16.555	2247	2249	2250	rBV2	2365	1434	0.14%	0.018%
81	16.684	2269	2271	2272	rVB2	3022	1627	0.16%	0.021%
82	16.737	2277	2280	2281	rBV3	4400	3555	0.35%	0.046%
83	16.896	2305	2307	2308	rBV2	2437	2064	0.20%	0.027%
84	17.037	2327	2331	2336	rBV	40152	52167	5.07%	0.672%
85	17.178	2353	2355	2357	rBV3	5536	5603	0.54%	0.072%
86	17.436	2393	2399	2404	rBV	562675	834394	81.11%	10.741%
87	17.483	2404	2407	2412	rVB	42757	59012	5.74%	0.760%
88	19.487	2744	2748	2753	rVB	97566	134562	13.08%	1.732%
89	19.734	2786	2790	2796	rVB	165989	182155	17.71%	2.345%
90	19.816	2801	2804	2806	rVV	52919	59651	5.80%	0.768%
91	19.845	2807	2809	2814	rVB2	61463	74677	7.26%	0.961%
92	20.339	2888	2893	2898	rVB4	73012	133634	12.99%	1.720%
93	20.774	2964	2967	2971	rVB	141206	156012	15.16%	2.008%
94	20.838	2975	2978	2982	rBV	112506	122579	11.92%	1.578%
95	21.350	3062	3065	3068	rVB	91473	101853	9.90%	1.311%
96	21.708	3121	3126	3139	rVB	549174	987656	96.00%	12.714%
97	21.808	3140	3143	3149	rVB	150841	185999	18.08%	2.394%
98	21.890	3152	3157	3161	rBV2	202316	415154	40.35%	5.344%
99	21.996	3171	3175	3180	rVB	257847	377243	36.67%	4.856%
100	24.951	3672	3678	3690	rVB2	341079	1028765	100.00%	13.243%

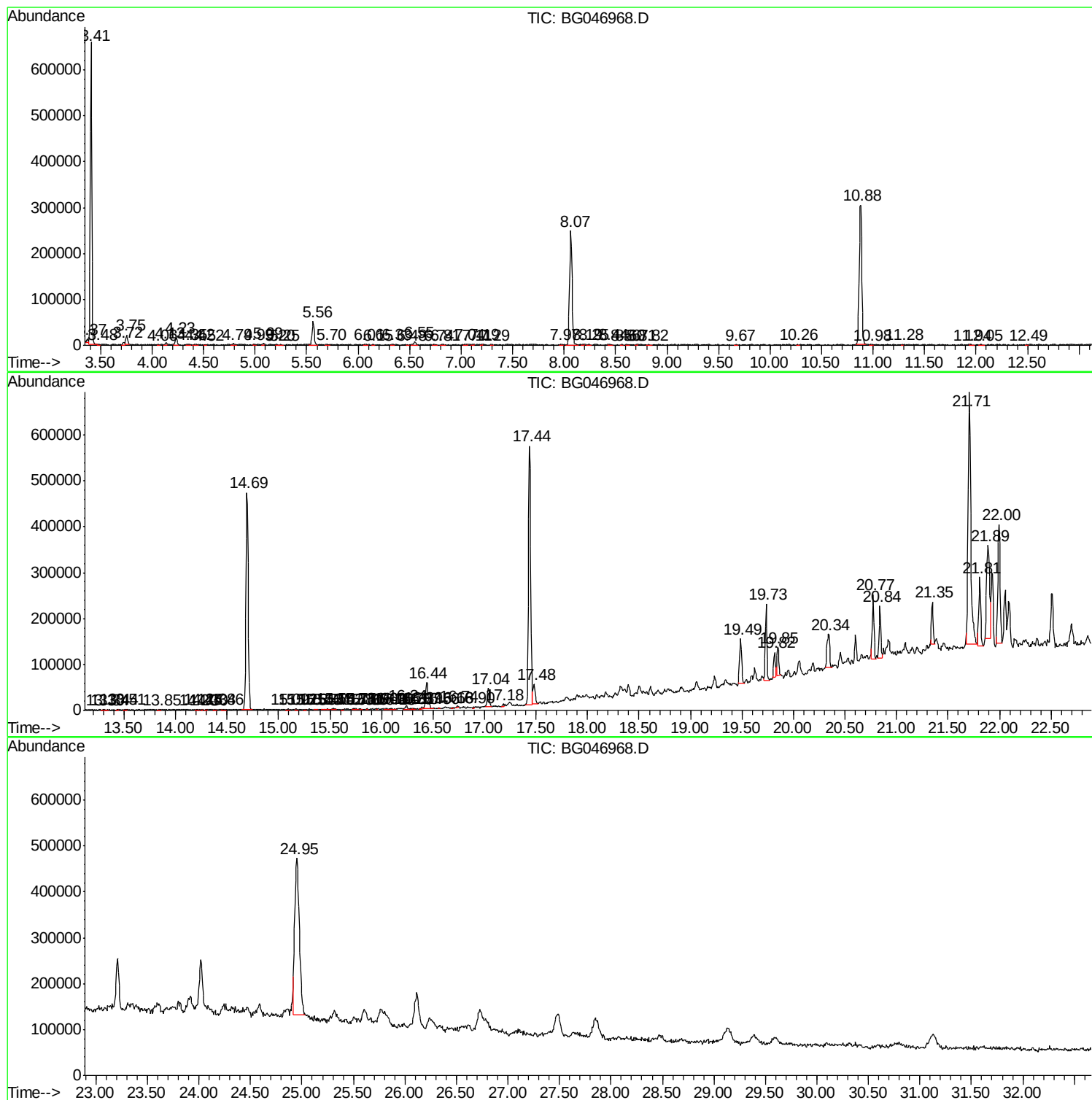
Sum of corrected areas: 7768310

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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG101520MA.M
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P



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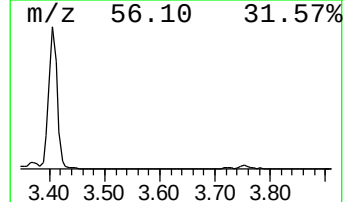
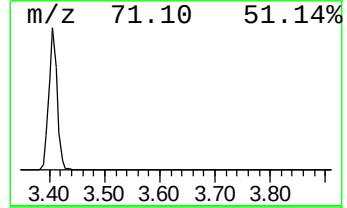
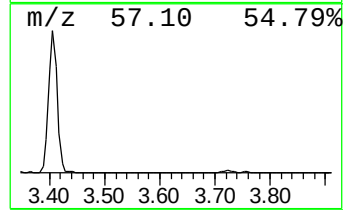
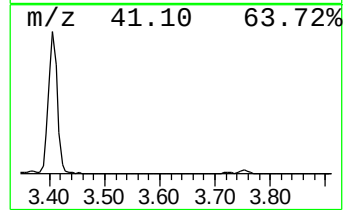
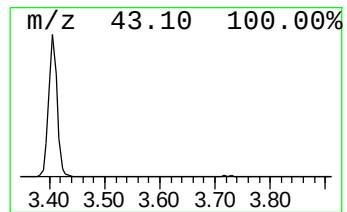
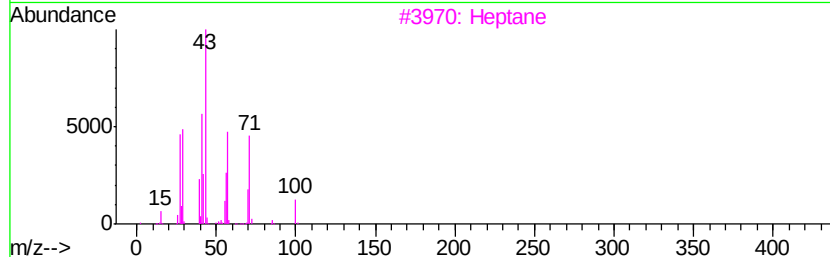
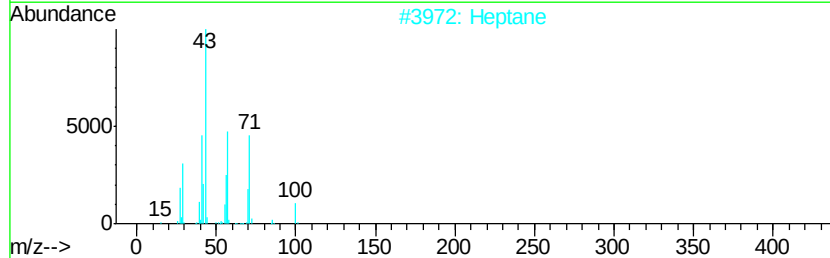
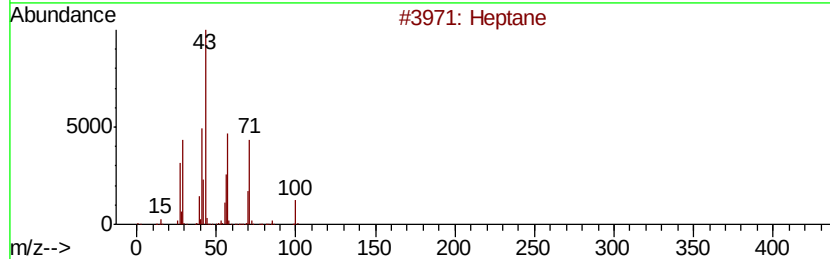
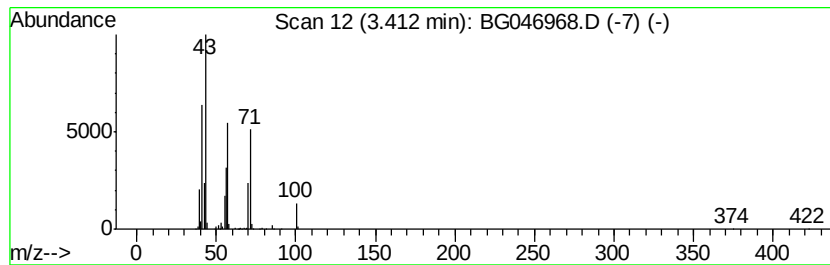
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG101520MA.M
Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.41	34.84 ng/ul	721177	1,4-Dichlorobenzene-d4	8.07

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heptane	100	C7H16	000142-82-5	91
2		Heptane	100	C7H16	000142-82-5	91
3		Heptane	100	C7H16	000142-82-5	91
4		Heptane	100	C7H16	000142-82-5	80
5		Oxalic acid, isobutyl pentyl ester	216	C11H20O4	1000309-37-0	45



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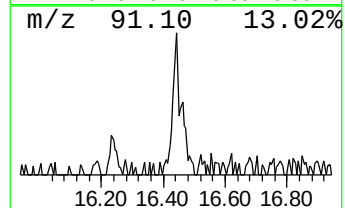
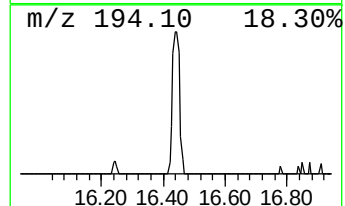
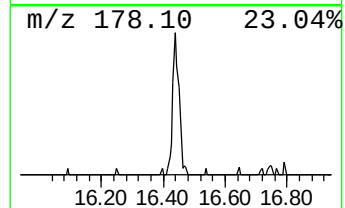
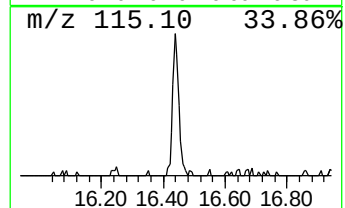
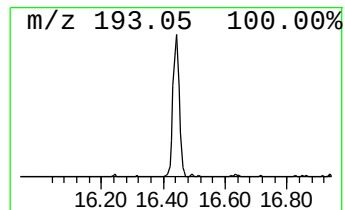
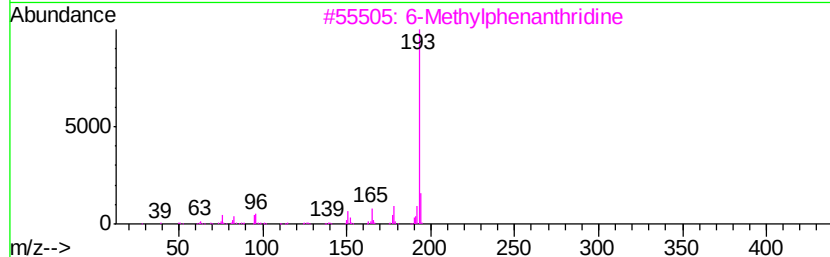
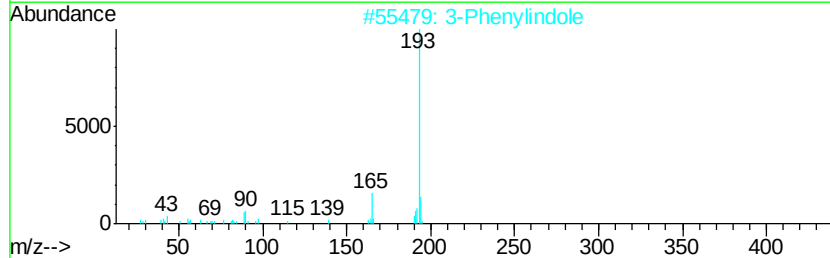
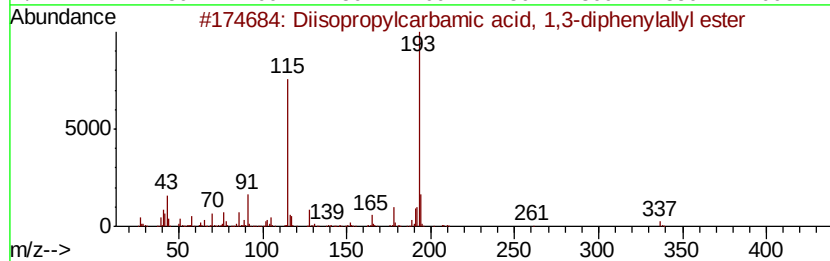
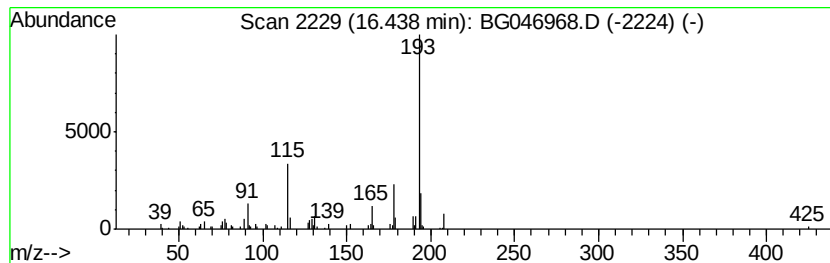
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 3 Diisopropylcarbamic acid, 1... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.44	2.35 ng/ul	97915	Phenanthrene-d10	17.44

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Diisopropylcarbamic acid, 1,3-di...	337	C22H27NO2	1000192-89-6	59
2		3-Phenylindole	193	C14H11N	001504-16-1	58
3		6-Methylphenanthridine	193	C14H11N	003955-65-5	53
4		1H-Indole, 2-phenyl-	193	C14H11N	000948-65-2	52
5		2-Aminophenanthrene	193	C14H11N	003366-65-2	50



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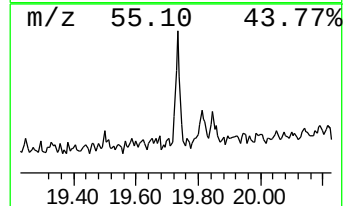
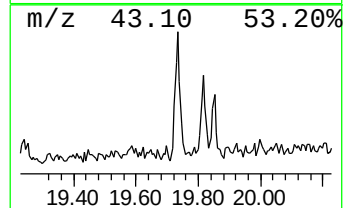
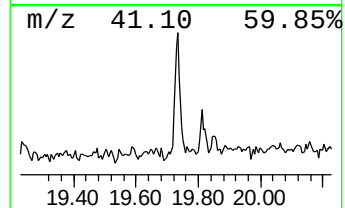
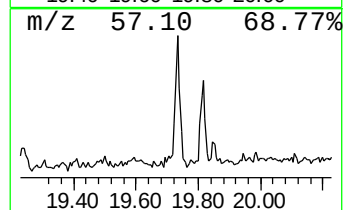
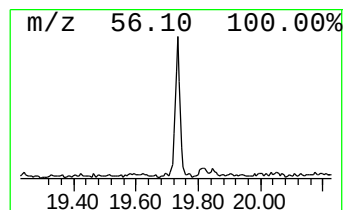
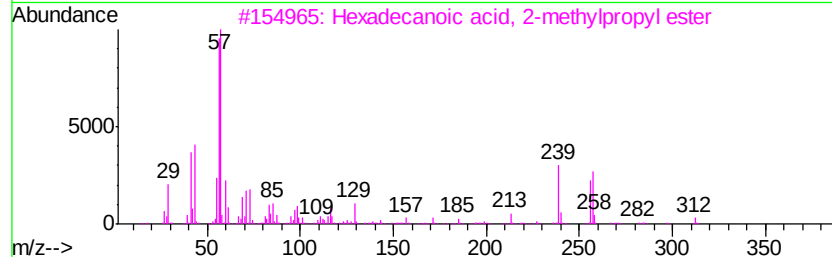
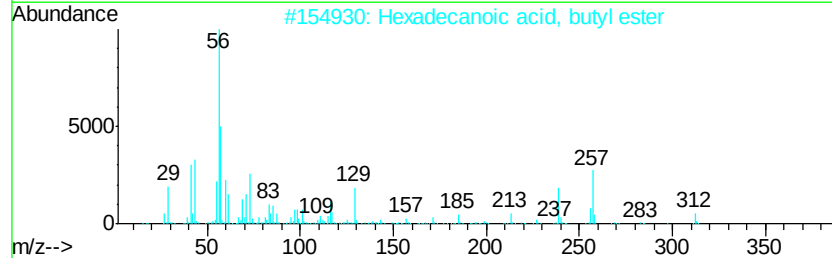
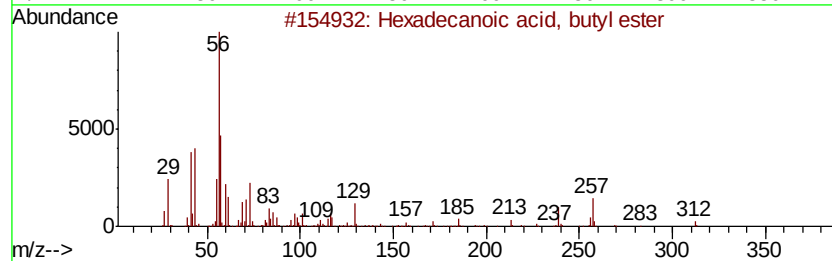
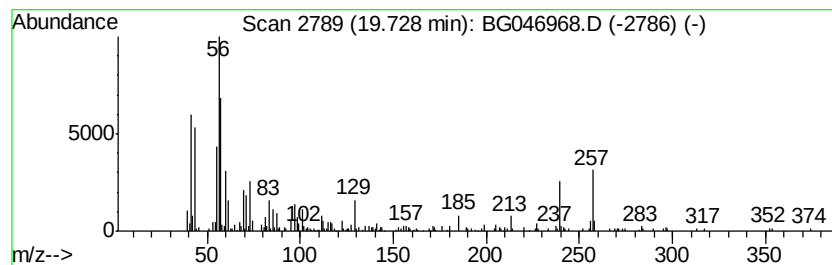
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Peak Number 4 Hexadecanoic acid, butyl ester Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.73	3.69 ng/ul	182155	Chrysene-d12	21.71

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexadecanoic acid, butyl ester	312	C20H40O2	000111-06-8	93
2		Hexadecanoic acid, butyl ester	312	C20H40O2	000111-06-8	91
3		Hexadecanoic acid, 2-methylpropyl...	312	C20H40O2	000110-34-9	89
4		Hexadecanoic acid, 1,1-dimethyle...	312	C20H40O2	031158-91-5	83
5		Eicosanoic acid	312	C20H40O2	000506-30-9	53



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG101720\
Data File : BG046968.D
Acq On : 18 Oct 2020 10:26
Operator : CG/JU
Sample : L4201-13
Misc : GCMS CONFIRMATION
ALS Vial : 34 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
C0AP3

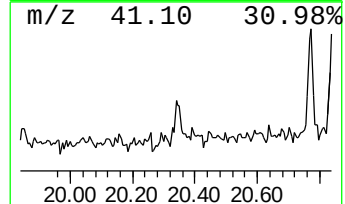
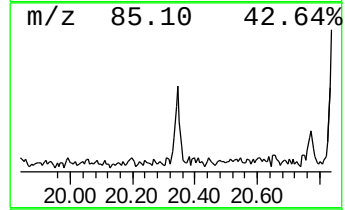
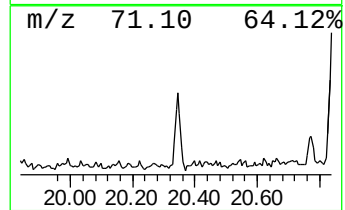
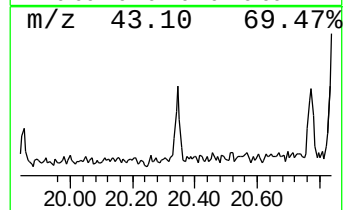
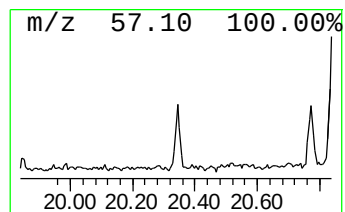
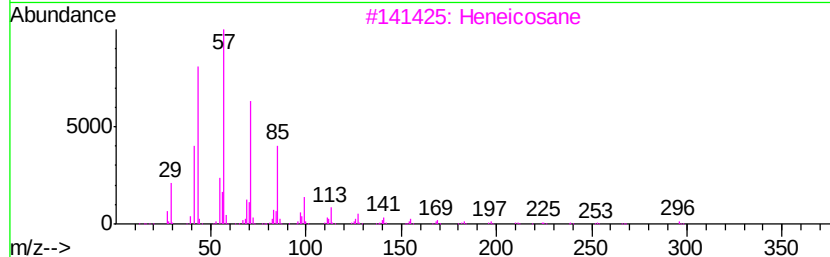
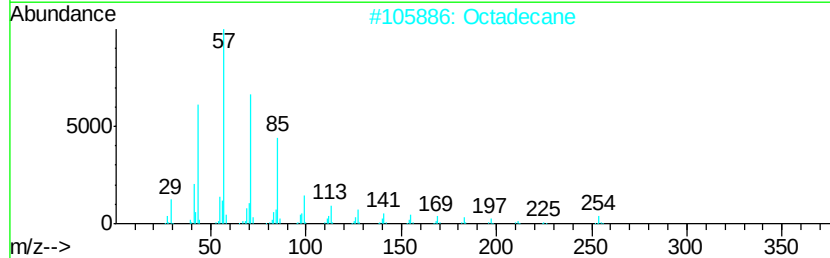
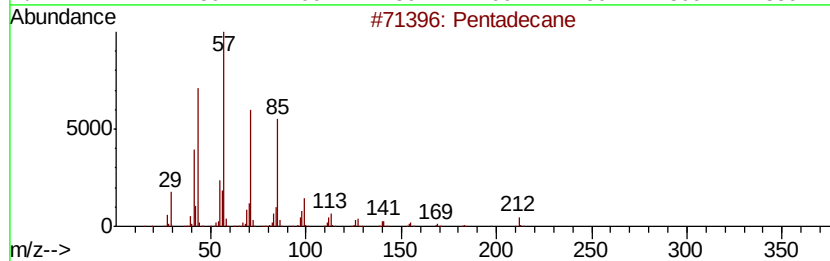
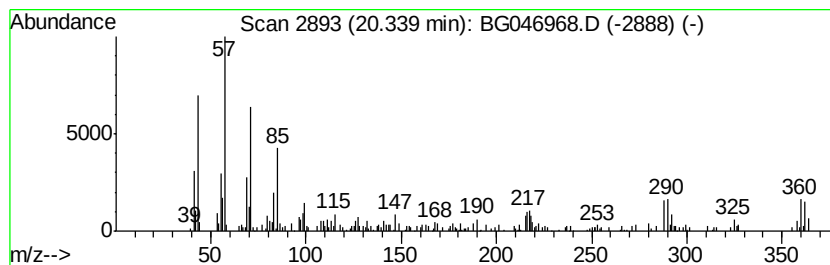
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 (DEL) Alkane: Straight-Chai... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.34	2.71 ng/ul	133634	Chrysene-d12	21.71

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Pentadecane	212	C15H32	000629-62-9	70
2		Octadecane	254	C18H38	000593-45-3	70
3		Heneicosane	296	C21H44	000629-94-7	70
4		Decane, 1-iodo-	268	C10H21I	002050-77-3	55
5		Tridecane	184	C13H28	000629-50-5	55



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG101720\
Data File : BG046968.D
Acq On : 18 Oct 2020 10:26
Operator : CG/JU
Sample : L4201-13
Misc : GCMS CONFIRMATION
ALS Vial : 34 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
C0AP3

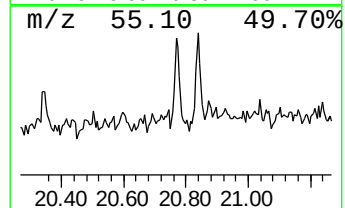
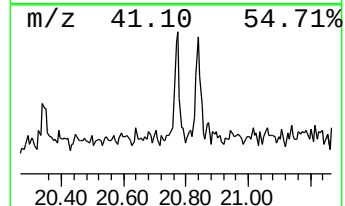
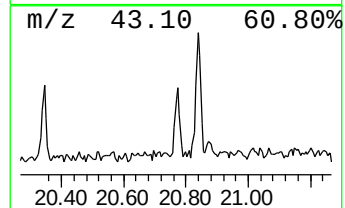
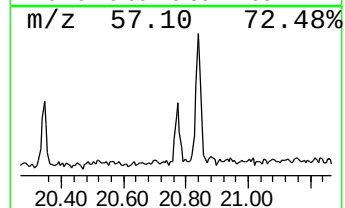
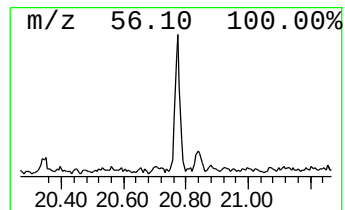
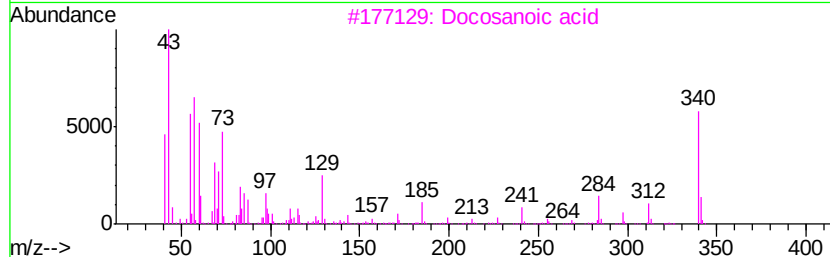
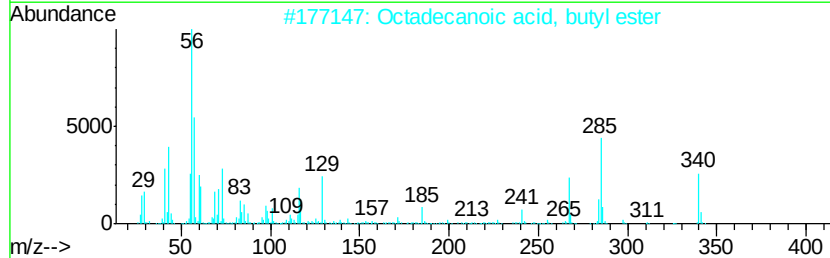
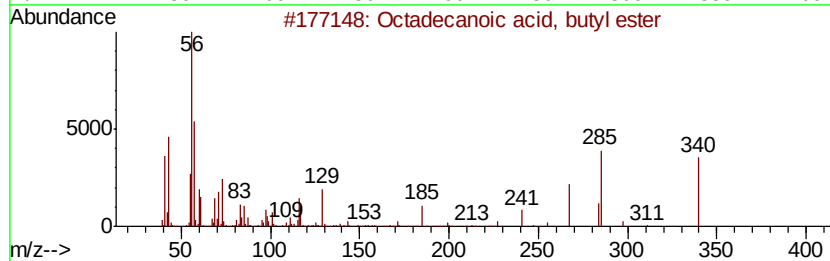
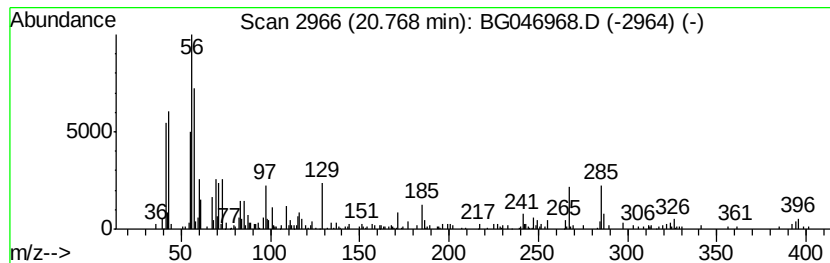
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 Octadecanoic acid, butyl ester Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.77	3.16 ng/ul	156012	Chrysene-d12	21.71

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octadecanoic acid, butyl ester	340	C22H44O2	000123-95-5	80
2		Octadecanoic acid, butyl ester	340	C22H44O2	000123-95-5	72
3		Docosanoic acid	340	C22H44O2	000112-85-6	55
4		Octadecanoic acid, 2-methylpropyl...	340	C22H44O2	000646-13-9	46
5		Butyl 15-methylhexadecanoate	326	C21H42O2	1000336-52-4	37



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG101720\
Data File : BG046968.D
Acq On : 18 Oct 2020 10:26
Operator : CG/JU
Sample : L4201-13
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ALS Vial : 34 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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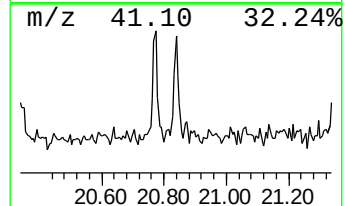
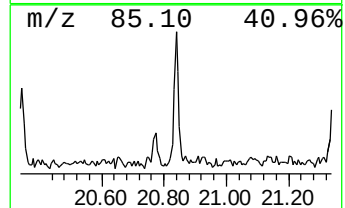
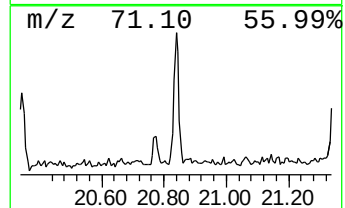
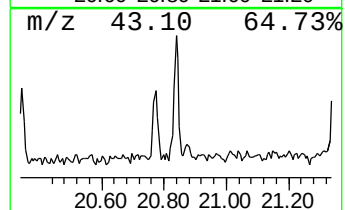
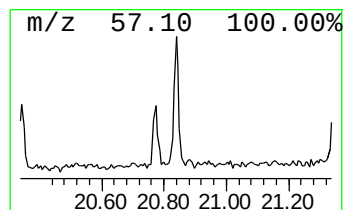
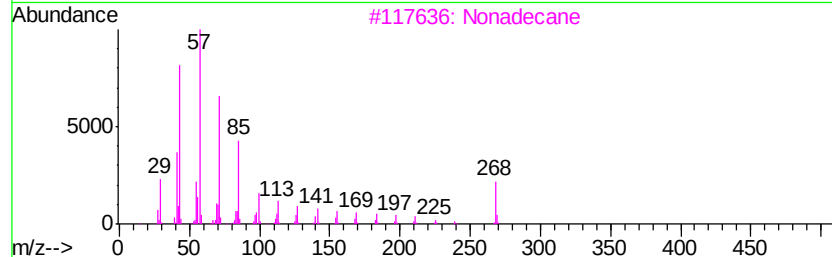
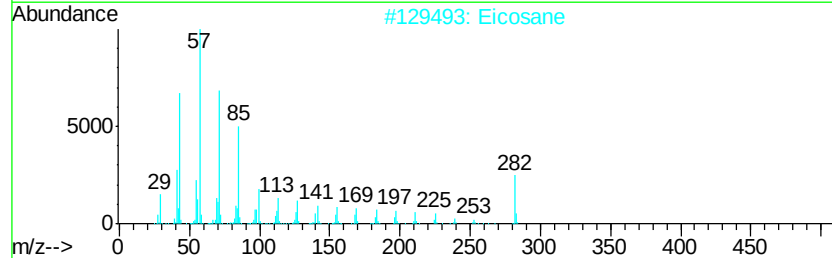
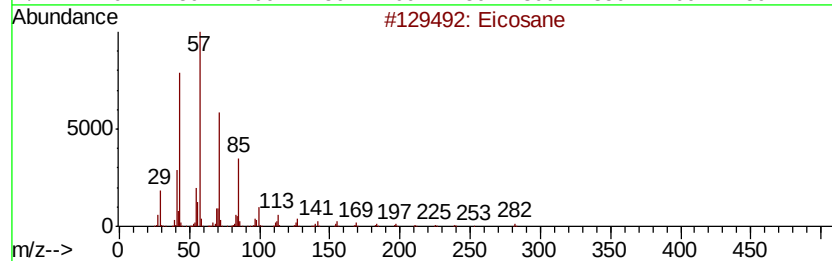
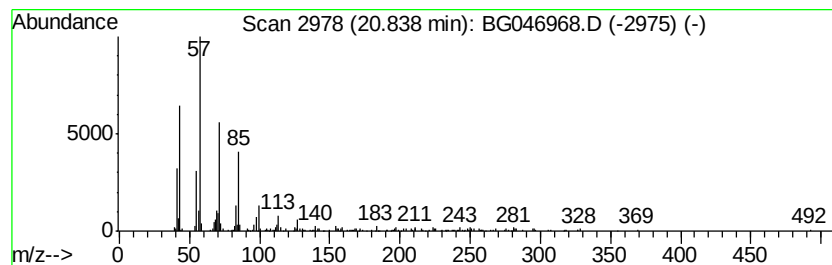
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 (DEL) Alkane: Straight-Chai... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.84	2.48 ng/ul	122579	Chrysene-d12	21.71

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Eicosane	282	C20H42	000112-95-8	87
2		Eicosane	282	C20H42	000112-95-8	87
3		Nonadecane	268	C19H40	000629-92-5	83
4		Nonadecane	268	C19H40	000629-92-5	80
5		Dodecane, 1-iodo-	296	C12H25I	004292-19-7	80



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG101720\
Data File : BG046968.D
Acq On : 18 Oct 2020 10:26
Operator : CG/JU
Sample : L4201-13
Misc : GCMS CONFIRMATION
ALS Vial : 34 Sample Multiplier: 1

Instrument :
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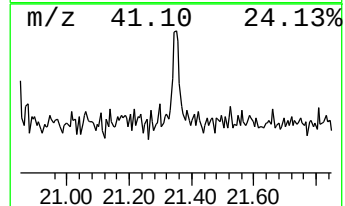
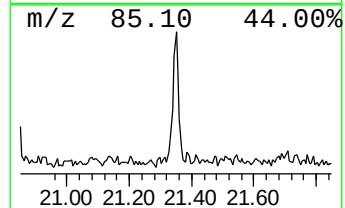
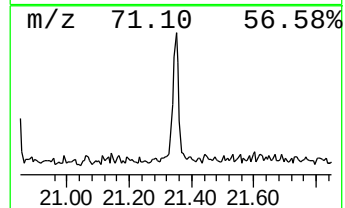
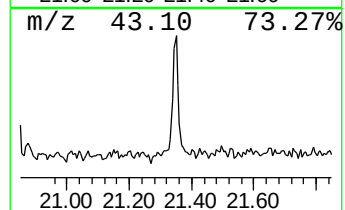
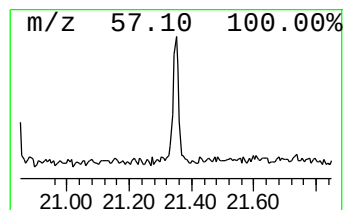
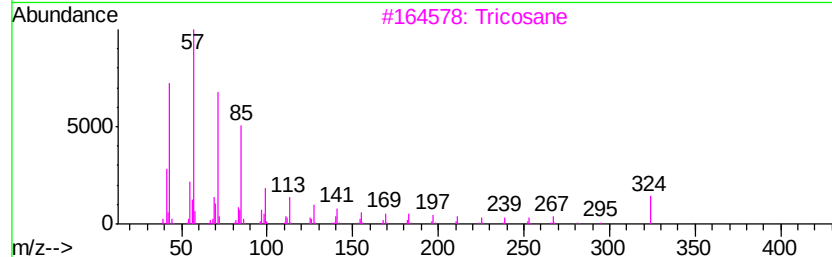
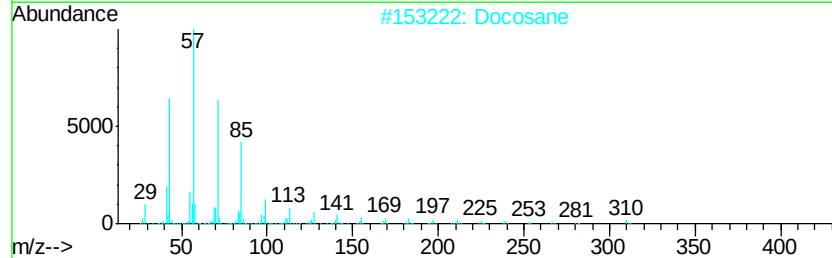
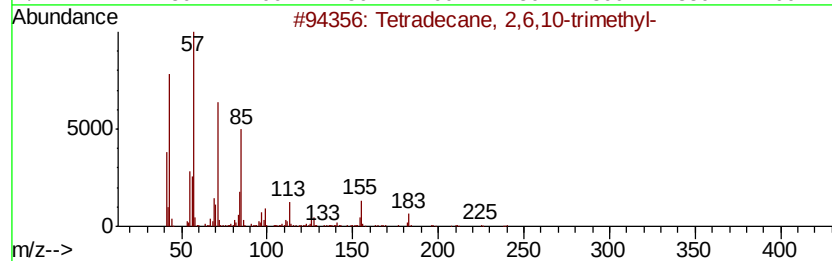
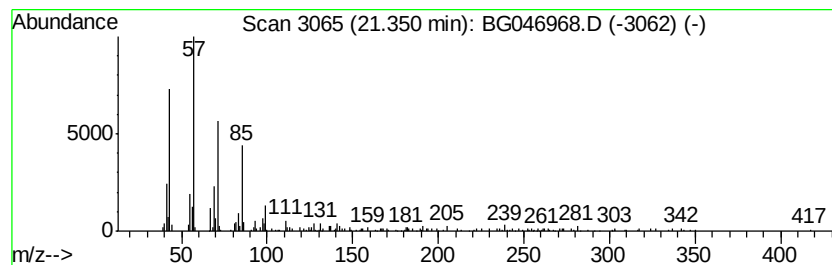
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 (DEL) Alkane: Straight-Chai... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.35	2.06 ng/ul	101853	Chrysene-d12	21.71

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tetradecane, 2,6,10-trimethyl-	240	C17H36	014905-56-7	87
2		Docosane	310	C22H46	000629-97-0	87
3		Tricosane	324	C23H48	000638-67-5	86
4		Tridecane, 1-iodo-	310	C13H27I	035599-77-0	80
5		Heptadecane	240	C17H36	000629-78-7	80



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG101720\
Data File : BG046968.D
Acq On : 18 Oct 2020 10:26
Operator : CG/JU
Sample : L4201-13
Misc : GCMS CONFIRMATION
ALS Vial : 34 Sample Multiplier: 1

Instrument :
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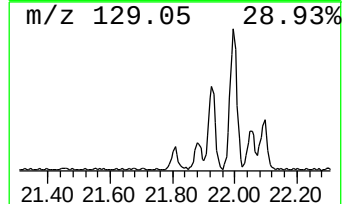
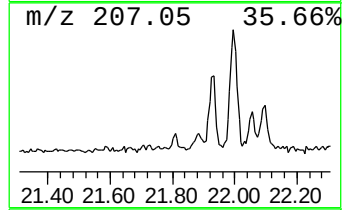
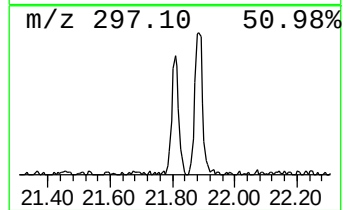
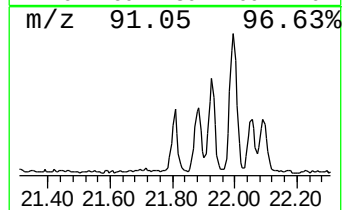
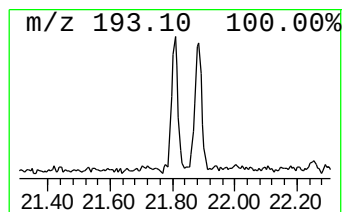
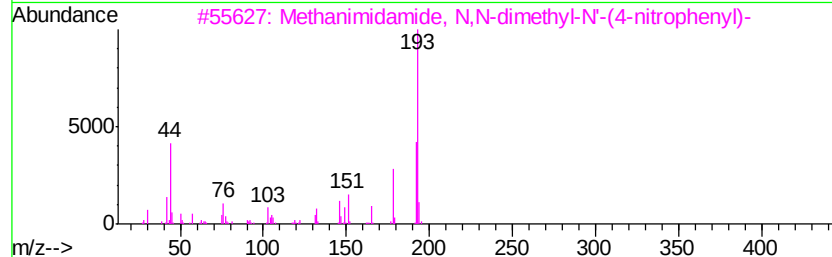
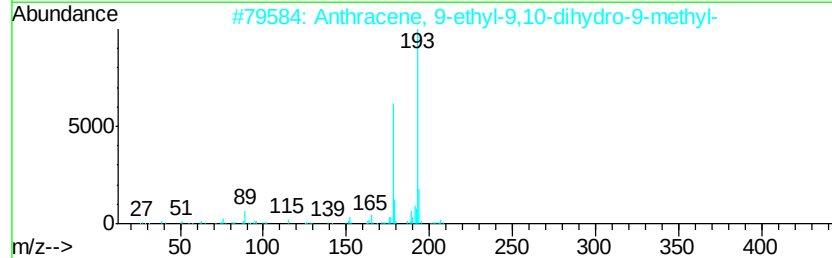
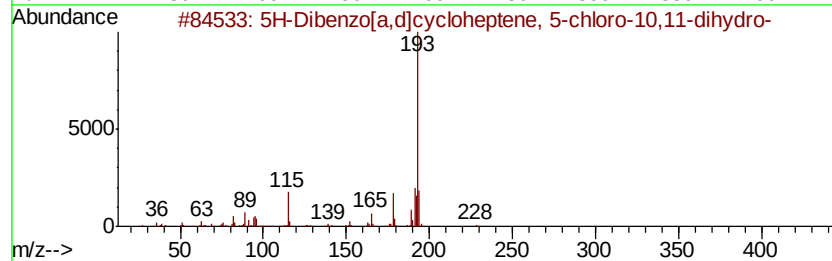
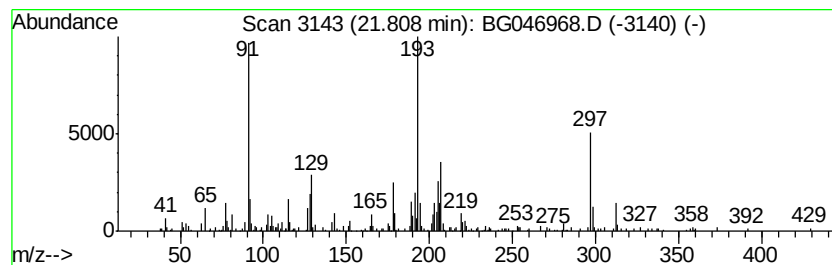
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 unknown-01 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.81	3.77 ng/ul	185999	Chrysene-d12	21.71

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	5H-Dibenzo[a,d]cycloheptene, 5-c...	228	C15H13Cl	001210-33-9	38
2		Anthracene, 9-ethyl-9,10-dihydro...	222	C17H18	054947-85-2	35
3		Methanimidamide, N,N-dimethyl-N'...	193	C9H11N3O2	001205-59-0	35
4		1-Phenylmethyl-2,4-diphenylimida...	312	C22H20N2	1000216-06-0	30
5		1,3-Dimethyl-6,8-dinitro-4-azaph...	297	C15H11N3O4	1000298-83-2	30



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG101720\
Data File : BG046968.D
Acq On : 18 Oct 2020 10:26
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Instrument :
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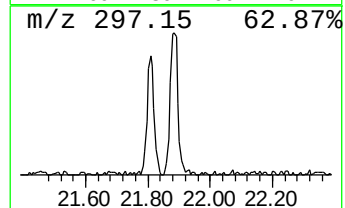
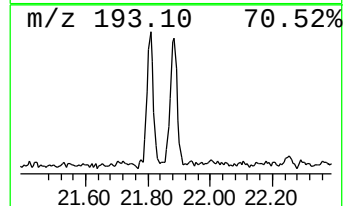
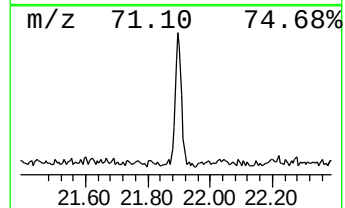
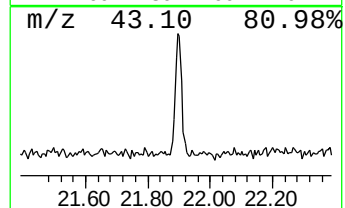
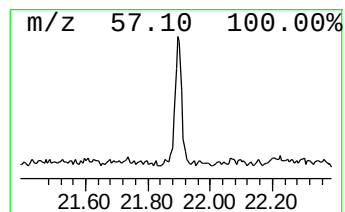
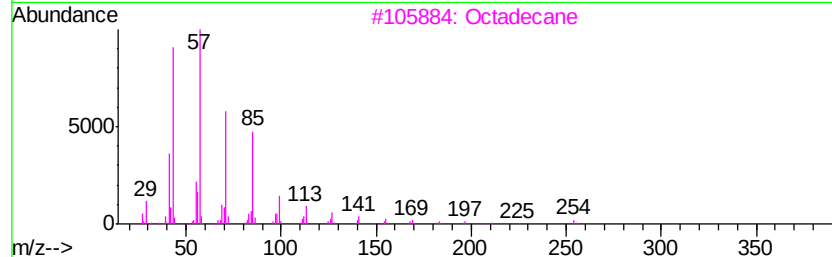
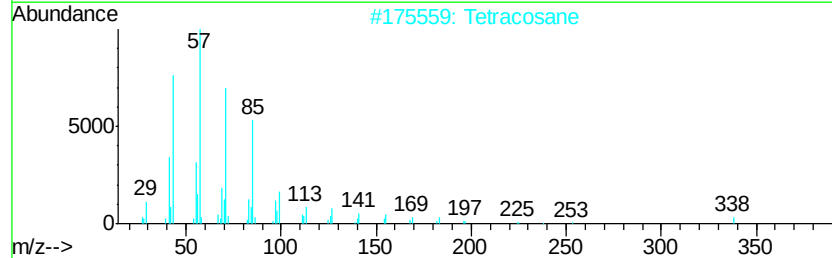
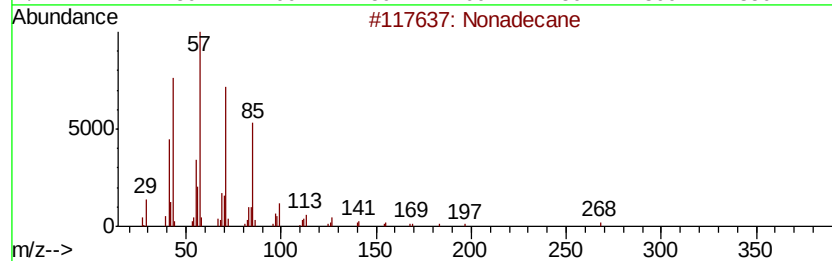
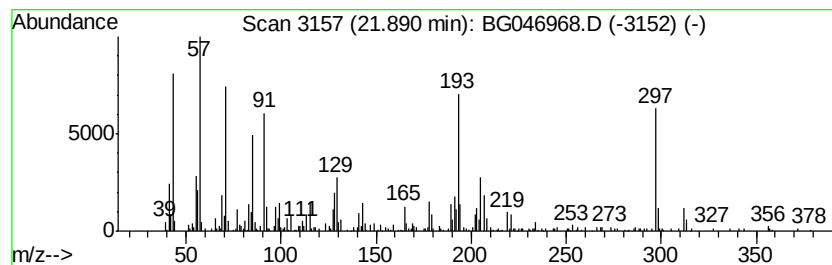
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 (DEL) Alkane: Straight-Chai... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.89	8.41 ng/ul	415154	Chrysene-d12	21.71

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Nonadecane	268	C19H40	000629-92-5	55
2		Tetracosane	338	C24H50	000646-31-1	42
3		Octadecane	254	C18H38	000593-45-3	38
4		Undecane, 5-methyl-	170	C12H26	001632-70-8	38
5		Docosane, 11-decyl-	451	C32H66	055401-55-3	38



Data Path : Z:\SVOASRV\HPCHEM1\BNA_G\DATA\BG101720\
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 ALS Vial : 34 Sample Multiplier: 1

Instrument :
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 ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG101520MA.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: Str...	3.41	34.8	ng/ul	721177	1	8.07	413985	20.0
Diisopropylcarbam...	16.44	2.4	ng/ul	97915	4	17.44	834394	20.0
Hexadecanoic acid...	19.73	3.7	ng/ul	182155	5	21.71	987656	20.0
(DEL) Alkane: Str...	20.34	2.7	ng/ul	133634	5	21.71	987656	20.0
Octadecanoic acid...	20.77	3.2	ng/ul	156012	5	21.71	987656	20.0
(DEL) Alkane: Str...	20.84	2.5	ng/ul	122579	5	21.71	987656	20.0
(DEL) Alkane: Str...	21.35	2.1	ng/ul	101853	5	21.71	987656	20.0
unknown-01	21.81	3.8	ng/ul	185999	5	21.71	987656	20.0
(DEL) Alkane: Str...	21.89	8.4	ng/ul	415154	5	21.71	987656	20.0
Methadone N-oxide	22.00	7.6	ng/ul	377243	5	21.71	987656	20.0