

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM061720\
 Data File : BM026435.D
 Acq On : 17 Jun 2020 20:43
 Operator : JU/CG
 Sample : L2959-08
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 BG2K2

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_M\METHODS\SOM-EPA-BM061220MA.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	3.669	128	133	144	rVB	112603	163664	0.20%	0.090%
2	6.540	617	621	624	rVV2	56127	93058	0.12%	0.051%
3	6.598	625	631	641	rVB	308536	550560	0.68%	0.301%
4	6.745	650	656	666	rBV	272920	472032	0.59%	0.258%
5	6.934	683	688	701	rVB	330531	590406	0.73%	0.323%
6	7.040	701	706	716	rBV2	122434	210889	0.26%	0.115%
7	7.392	758	766	774	rVB	614533	957446	1.19%	0.524%
8	7.616	796	804	809	rBV2	208269	339498	0.42%	0.186%
9	7.669	809	813	822	rVB5	43717	94301	0.12%	0.052%
10	7.934	854	858	864	rVB2	69088	104944	0.13%	0.057%
11	8.116	884	889	895	rVB	412361	621632	0.77%	0.340%
12	8.492	940	953	956	rBV9	33383	113249	0.14%	0.062%
13	8.539	956	961	967	rVV	348823	573519	0.71%	0.314%
14	8.628	971	976	983	rVV	151871	235276	0.29%	0.129%
15	8.733	983	994	1001	rVB9	38351	119871	0.15%	0.066%
16	9.251	1072	1082	1088	rBV	305556	512226	0.64%	0.280%
17	9.539	1126	1131	1139	rBV3	50314	102680	0.13%	0.056%
18	9.792	1165	1174	1182	rBV	419849	715049	0.89%	0.391%
19	10.151	1228	1235	1244	rVB	864553	1393490	1.73%	0.762%
20	10.304	1254	1261	1276	rBV	313625	627377	0.78%	0.343%
21	10.704	1324	1329	1340	rBV	77263	166571	0.21%	0.091%
22	12.227	1583	1588	1593	rBV2	47487	88275	0.11%	0.048%
23	12.280	1593	1597	1605	rVB	51861	81693	0.10%	0.045%
24	12.539	1635	1641	1648	rBV	286997	466022	0.58%	0.255%
25	12.774	1676	1681	1690	rBV	117136	197340	0.25%	0.108%
26	13.086	1726	1734	1743	rBV	173335	279462	0.35%	0.153%
27	13.486	1795	1802	1806	rBV	866883	1264379	1.57%	0.691%
28	13.527	1806	1809	1820	rVB	239586	367099	0.46%	0.201%
29	13.733	1838	1844	1855	rBV	894491	1333510	1.66%	0.729%
30	14.051	1890	1898	1909	rVB	1318273	1932704	2.40%	1.057%
31	14.292	1933	1939	1942	rBV	367367	538173	0.67%	0.294%
32	14.333	1942	1946	1958	rVB	1234125	1782989	2.22%	0.975%
33	14.533	1975	1980	1989	rBV	350546	495340	0.62%	0.271%
34	14.610	1989	1993	1996	rVV	91260	120402	0.15%	0.066%

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Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM061220MA.M
 Title : SVOA CALIBRATION

35	14.651	1996	2000	2003	rVV	436712	652138	0.81%	0.357%
36	14.692	2003	2007	2014	rVB	2304624	3163940	3.93%	1.730%
37	15.051	2058	2068	2071	rBV	1122960	1693228	2.11%	0.926%
38	15.080	2071	2073	2079	rVB	1004640	1175631	1.46%	0.643%
39	15.221	2088	2097	2106	rBV	8051390	11992853	14.91%	6.558%
40	15.457	2131	2137	2145	rBV	612398	1008528	1.25%	0.552%
41	15.815	2194	2198	2200	rBV	91791	123742	0.15%	0.068%
42	15.839	2200	2202	2210	rVB	82527	107991	0.13%	0.059%
43	16.262	2270	2274	2280	rVB2	90550	124296	0.15%	0.068%
44	16.515	2303	2317	2321	rBV	30823501	80424910	100.00%	43.982%
45	16.609	2329	2333	2337	rVB	143237	191179	0.24%	0.105%
46	16.662	2338	2342	2350	rVB	96181	159600	0.20%	0.087%
47	16.815	2362	2368	2373	rBV	1713095	2297134	2.86%	1.256%
48	16.909	2379	2384	2393	rVB2	1266313	1692281	2.10%	0.925%
49	17.192	2427	2432	2437	rBV	313435	451018	0.56%	0.247%
50	17.345	2450	2458	2461	rBV2	217429	350710	0.44%	0.192%
51	17.427	2467	2472	2480	rVB3	123634	254264	0.32%	0.139%
52	17.762	2522	2529	2532	rBV	2012647	2807670	3.49%	1.535%
53	17.804	2532	2536	2542	rVB	3140047	3821956	4.75%	2.090%
54	18.039	2571	2576	2580	rBV	326609	396728	0.49%	0.217%
55	18.503	2652	2655	2662	rVB3	85969	132106	0.16%	0.072%
56	18.680	2681	2685	2690	rVB	640707	754187	0.94%	0.412%
57	18.933	2723	2728	2739	rBV	1192271	2983400	3.71%	1.632%
58	19.056	2745	2749	2755	rBV	1467645	1956054	2.43%	1.070%
59	19.115	2756	2759	2767	rVV	402366	637047	0.79%	0.348%
60	19.215	2771	2776	2782	rVB	1355820	1929944	2.40%	1.055%
61	19.268	2782	2785	2790	rVB	1024652	1154826	1.44%	0.632%
62	19.639	2845	2848	2851	rVB	168979	182419	0.23%	0.100%
63	19.809	2873	2877	2882	rVB	1325926	1404846	1.75%	0.768%
64	20.133	2929	2932	2935	rBV	93087	92510	0.12%	0.051%
65	20.309	2958	2962	2972	rVB	1593700	1721093	2.14%	0.941%
66	20.592	3006	3010	3011	rBV2	148689	194114	0.24%	0.106%
67	20.615	3011	3014	3019	rVB2	455450	559610	0.70%	0.306%
68	20.727	3029	3033	3037	rBV2	151467	286364	0.36%	0.157%
69	20.774	3037	3041	3046	rVB	2409954	2806571	3.49%	1.535%
70	20.968	3071	3074	3078	rBV3	193069	262845	0.33%	0.144%
71	21.021	3078	3083	3087	rVV	2304170	2701586	3.36%	1.477%

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72	21.215	3111	3116	3120	rBV	2125666	2289654	2.85%	1.252%
73	21.474	3158	3160	3163	rBV	149710	145160	0.18%	0.079%
74	21.633	3182	3187	3194	rBV	2545148	2910644	3.62%	1.592%
75	21.856	3221	3225	3230	rBV	312502	459956	0.57%	0.252%
76	22.062	3256	3260	3265	rVB	2945840	3337288	4.15%	1.825%
77	22.533	3335	3340	3352	rVB	2822595	3733498	4.64%	2.042%
78	22.980	3411	3416	3422	rBV	1143873	1893073	2.35%	1.035%
79	23.050	3422	3428	3433	rVV	2295593	3764405	4.68%	2.059%
80	23.115	3433	3439	3447	rVV	1803996	3149006	3.92%	1.722%
81	23.197	3449	3453	3457	rVV4	243514	405341	0.50%	0.222%
82	23.244	3458	3461	3479	rVB	214391	535864	0.67%	0.293%
83	23.633	3521	3527	3535	rBV	1740287	2984026	3.71%	1.632%
84	24.044	3593	3597	3608	rVB3	204086	452222	0.56%	0.247%
85	24.303	3635	3641	3652	rBV	919972	2025275	2.52%	1.108%
86	24.444	3660	3665	3674	rVB	428000	865723	1.08%	0.473%
87	24.550	3676	3683	3698	rVB2	686200	1773300	2.20%	0.970%
88	25.085	3768	3774	3780	rVV	418996	891005	1.11%	0.487%
89	25.162	3782	3787	3802	rVB	339694	890264	1.11%	0.487%

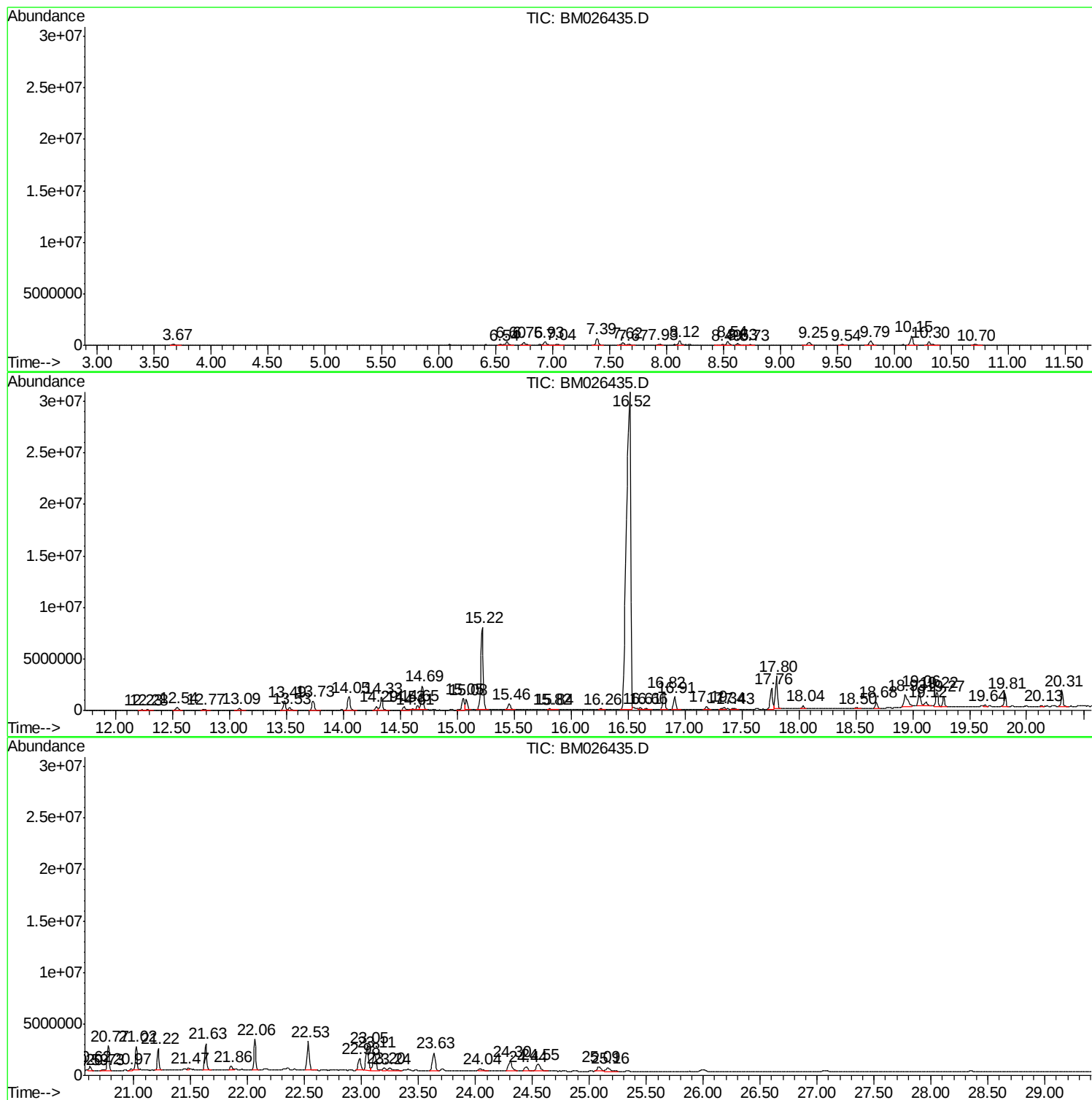
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM061220MA.M
Quant Title : SVOA CALIBRATION

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



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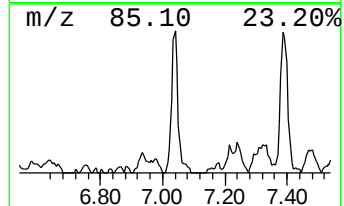
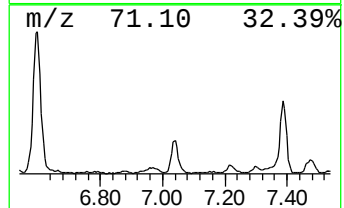
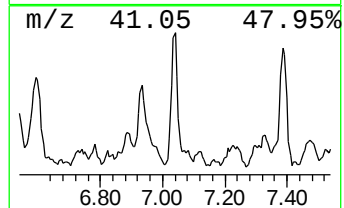
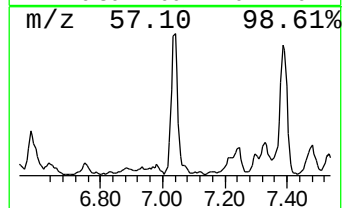
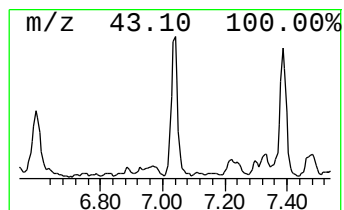
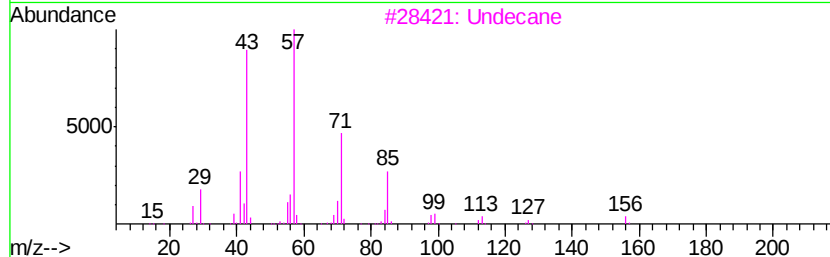
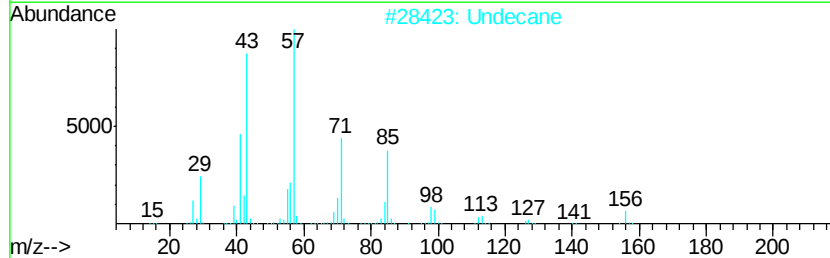
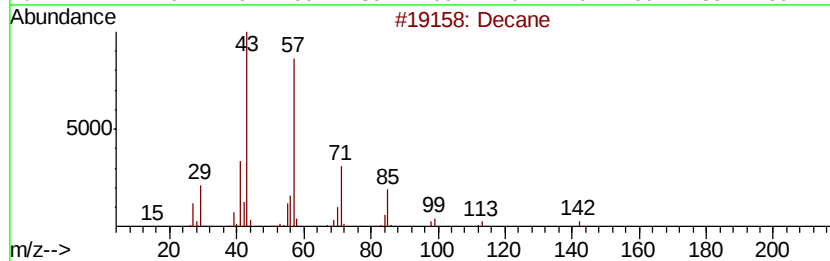
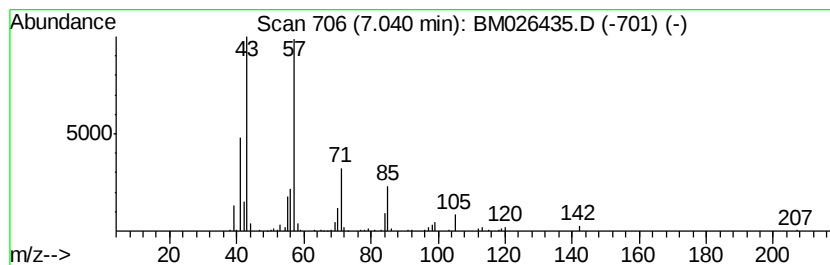
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM061220MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 (DEL) Alkane: Straight-Chai... Concentration Rank 24

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.04	4.41 ng/ul	210889	1,4-Dichlorobenzene-d4	7.39

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Decane	142	C10H22	000124-18-5	94
2		Undecane	156	C11H24	001120-21-4	83
3		Undecane	156	C11H24	001120-21-4	78
4		Nonane	128	C9H20	000111-84-2	72
5		Undecane	156	C11H24	001120-21-4	64



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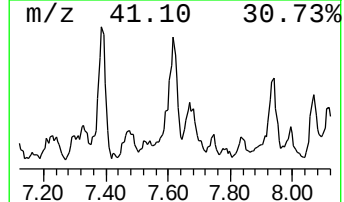
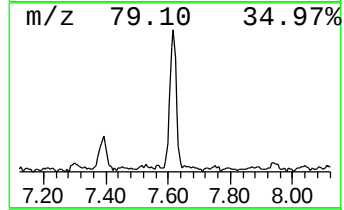
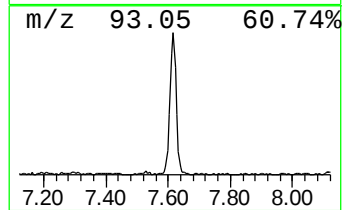
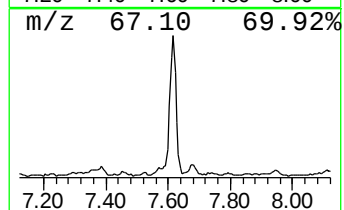
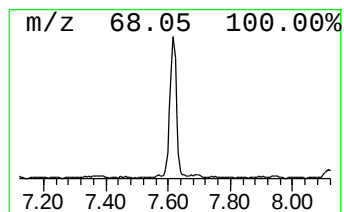
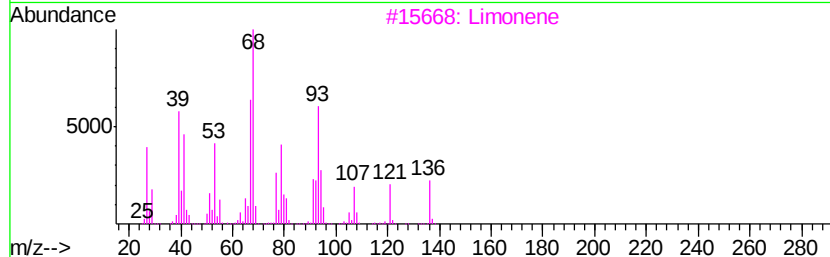
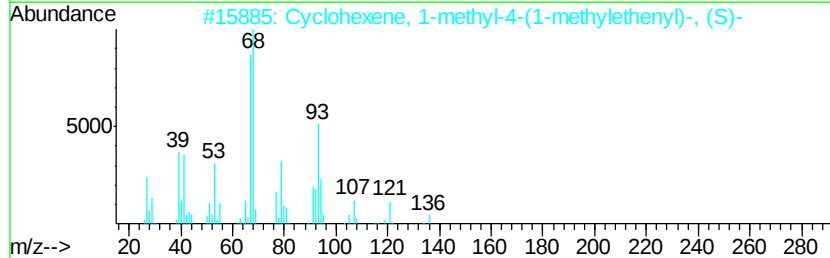
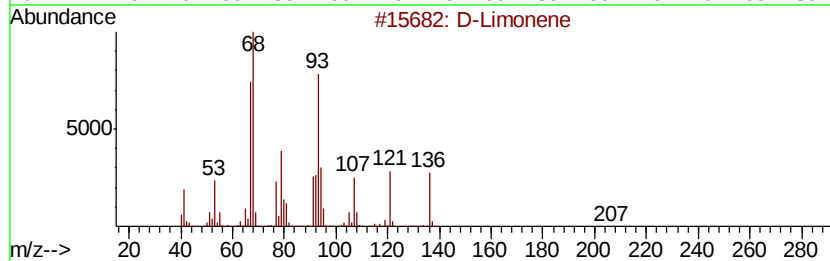
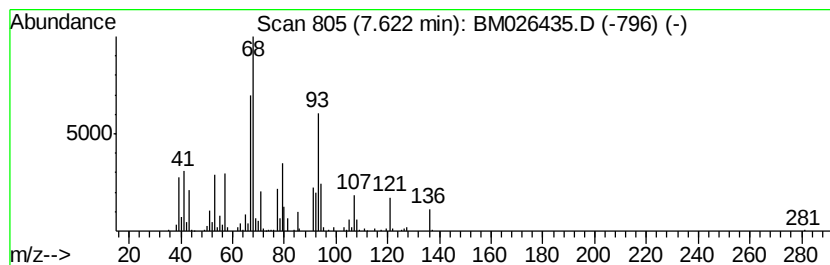
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM061220MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 D-Limonene Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.62	7.09 ng/ul	339498	1,4-Dichlorobenzene-d4	7.39

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	D-Limonene	136	C10H16	005989-27-5	95
2		Cyclohexene, 1-methyl-4-(1-methy...	136	C10H16	005989-54-8	94
3		Limonene	136	C10H16	000138-86-3	87
4		D-Limonene	136	C10H16	005989-27-5	83
5		Limonene	136	C10H16	000138-86-3	81



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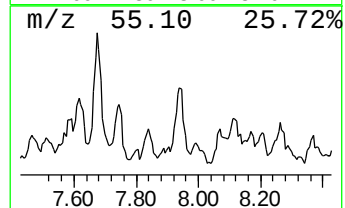
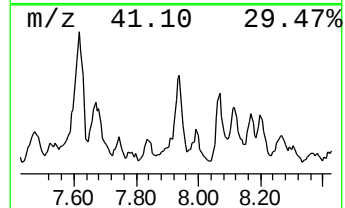
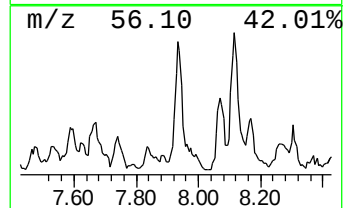
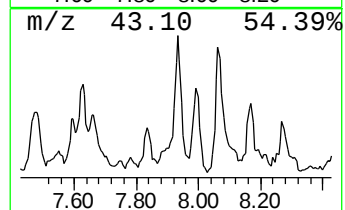
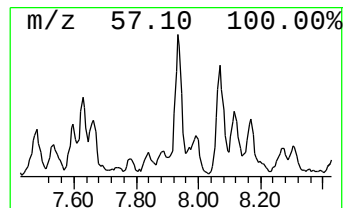
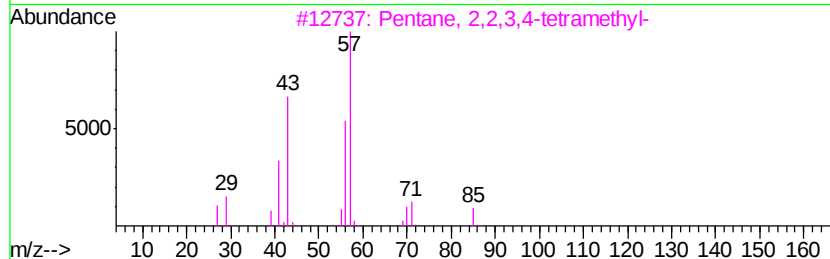
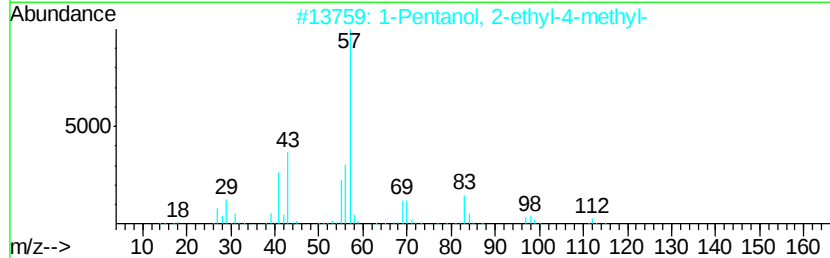
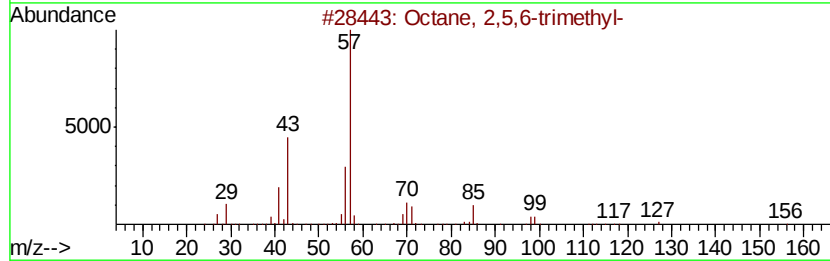
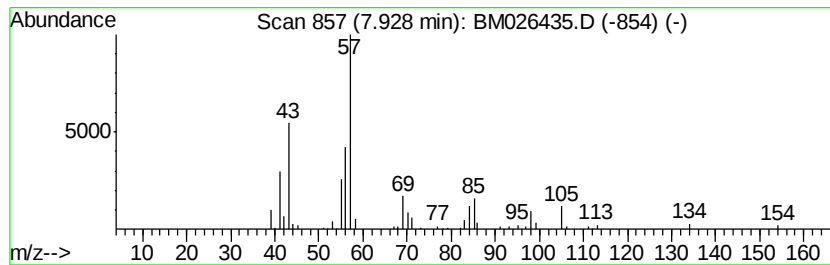
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM061220MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 4 (DEL) Alkane: Straight-Chai... Concentration Rank 38

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.93	2.19 ng/ul	104944	1,4-Dichlorobenzene-d4	7.39

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octane, 2,5,6-trimethyl-	156	C11H24	062016-14-2	59
2			1-Pentanol, 2-ethyl-4-methyl-	130	C8H18O	000106-67-2	50
3			Pentane, 2,2,3,4-tetramethyl-	128	C9H20	001186-53-4	50
4			Pentane, 2,2,3,4-tetramethyl-	128	C9H20	001186-53-4	50
5			Ether, heptyl hexyl	200	C13H28O	007289-40-9	50



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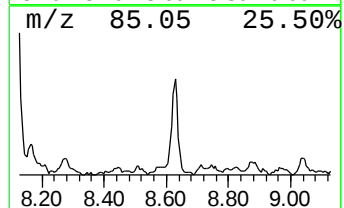
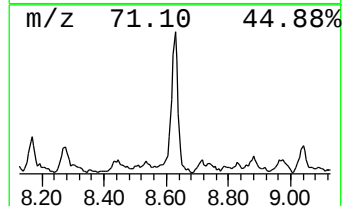
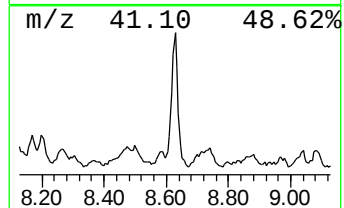
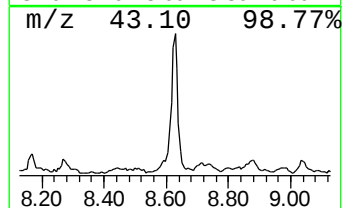
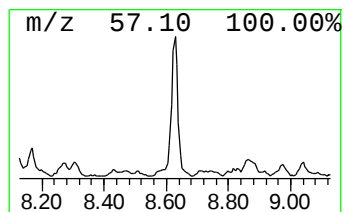
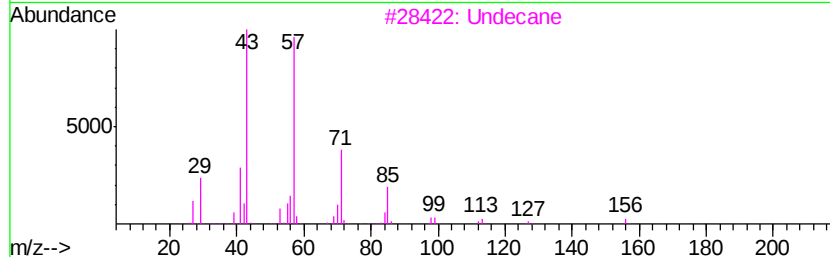
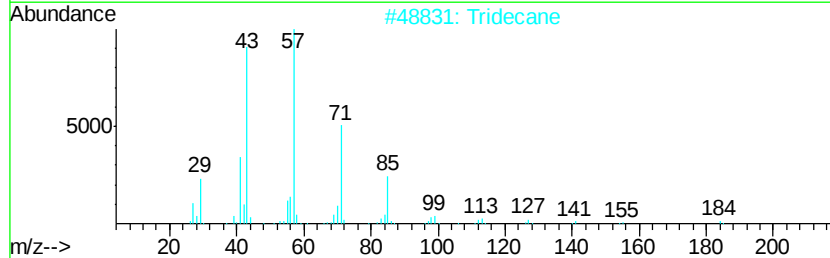
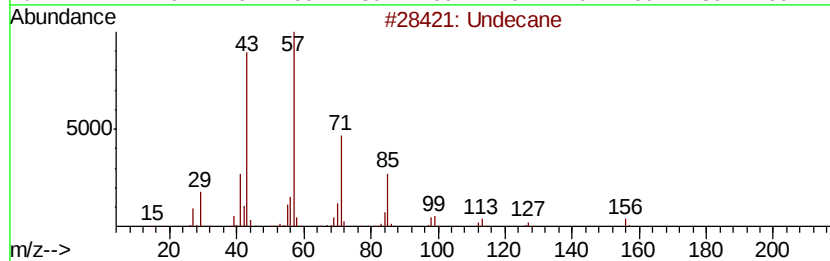
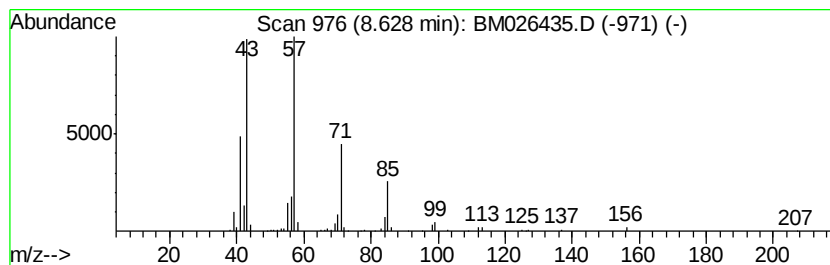
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 (DEL) Alkane: Straight-Chai... Concentration Rank 22

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.63	4.91 ng/ul	235276	1,4-Dichlorobenzene-d4	7.39

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Undecane			156	C11H24	001120-21-4	91
2	Tridecane			184	C13H28	000629-50-5	90
3	Undecane			156	C11H24	001120-21-4	90
4	Undecane			156	C11H24	001120-21-4	86
5	Tetradecane			198	C14H30	000629-59-4	83



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM061720\
Data File : BM026435.D
Acq On : 17 Jun 2020 20:43
Operator : JU/CG
Sample : L2959-08
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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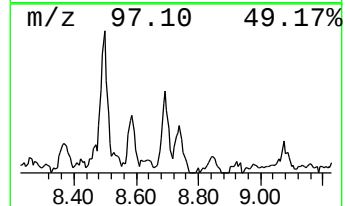
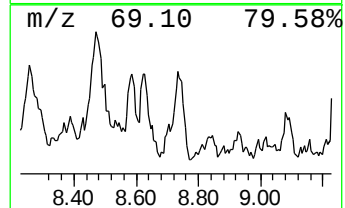
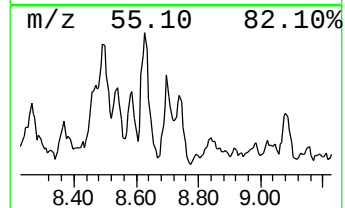
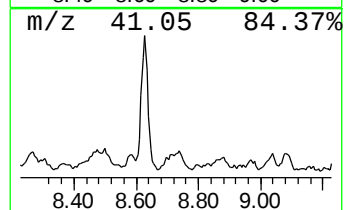
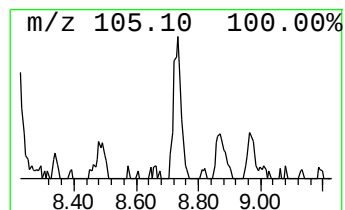
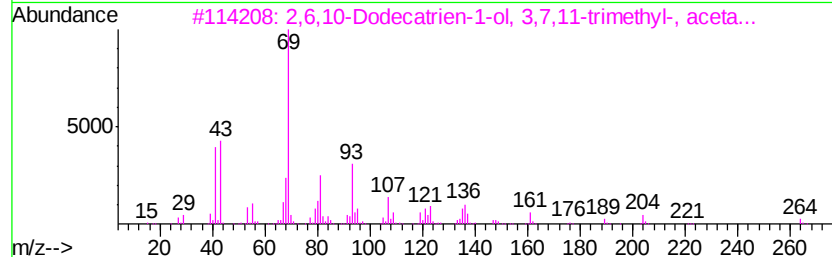
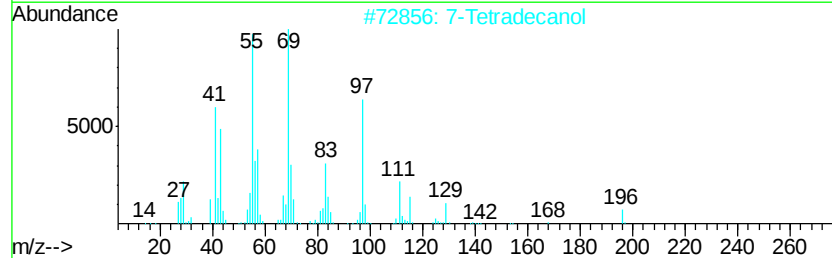
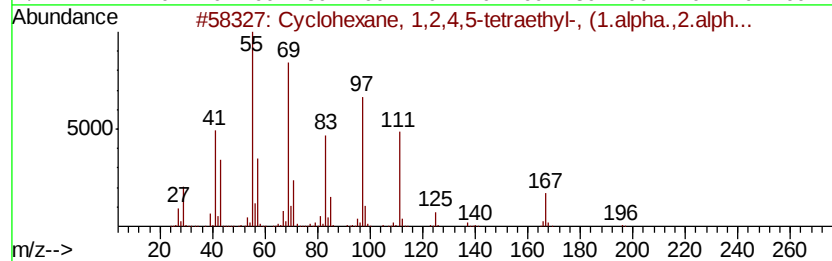
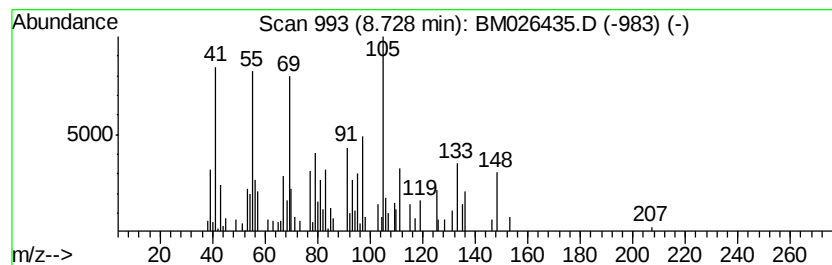
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 (DEL) Alkane: Cyclic8.73 Concentration Rank 35

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.73	2.50 ng/ul	119871	1,4-Dichlorobenzene-d4	7.39

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclohexane, 1,2,4,5-tetraethyl-...	196	C14H28	061142-24-3	38
2		7-Tetradecanol	214	C14H30O	003981-79-1	35
3		2,6,10-Dodecatrien-1-ol, 3,7,11-...	264	C17H28O2	004128-17-0	22
4		4-Hexen-1-ol, 5-methyl-2-(1-meth...	154	C10H18O	058461-27-1	22
5		Triallylsilane	152	C9H16Si	001116-62-7	22



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM061720\
Data File : BM026435.D
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Sample : L2959-08
Misc :
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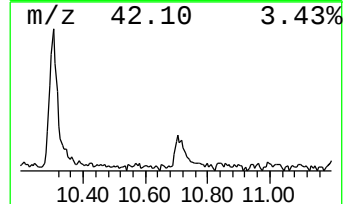
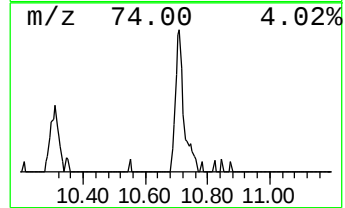
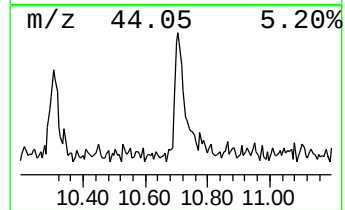
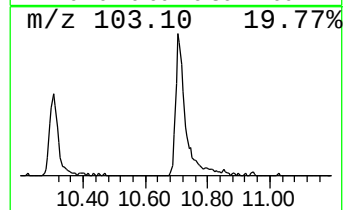
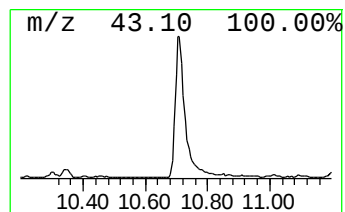
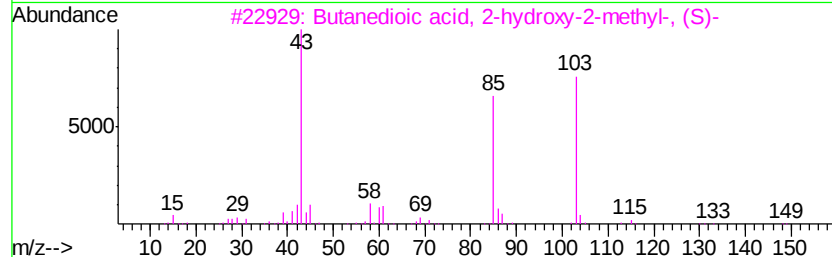
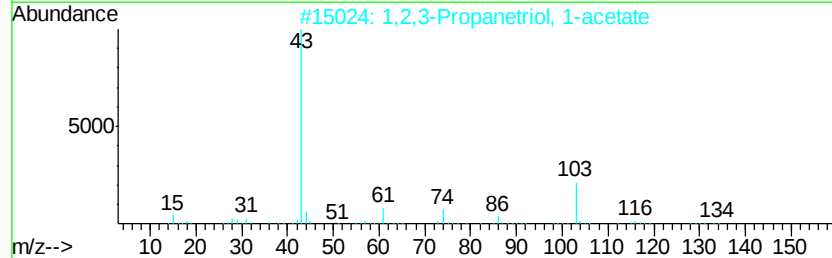
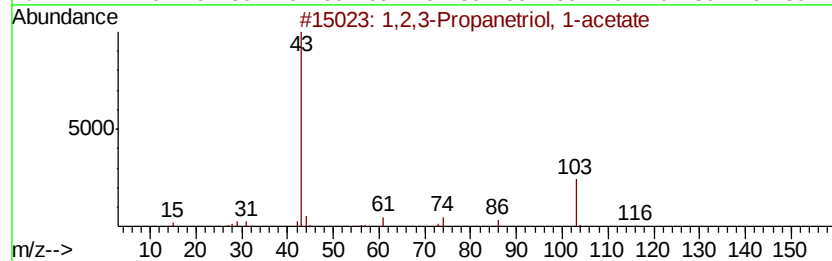
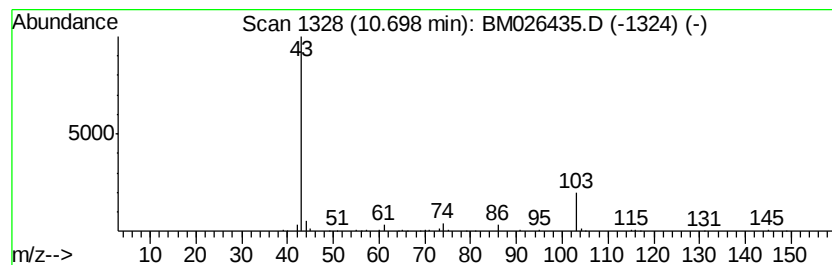
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 1,2,3-Propanetriol, 1-acetate Concentration Rank 36

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.70	2.39 ng/ul	166571	Naphthalene-d8	10.15

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1,2,3-Propanetriol, 1-acetate	134	C5H10O4	000106-61-6	50
2		1,2,3-Propanetriol, 1-acetate	134	C5H10O4	000106-61-6	38
3		Butanedioic acid, 2-hydroxy-2-methyl-, (S)-	148	C5H8O5	006236-09-5	9
4		2-Hydroxyimino-3-oxo-butanoic acid	145	C5H7NO4	073708-95-9	7
5		Ketone, methyl 2-methyl-1,3-oxot-	146	C6H10O2S	033266-06-7	4



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM061720\
Data File : BM026435.D
Acq On : 17 Jun 2020 20:43
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Misc :
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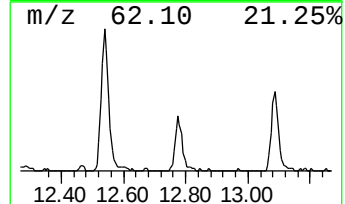
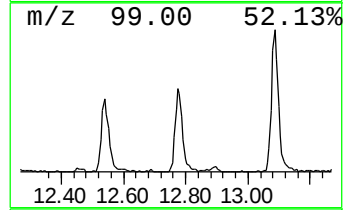
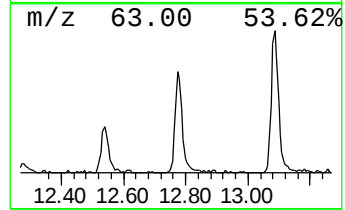
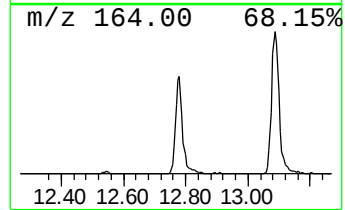
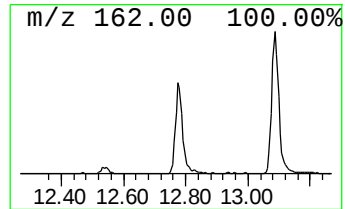
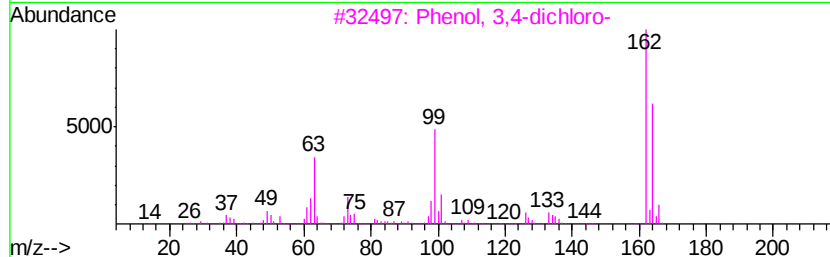
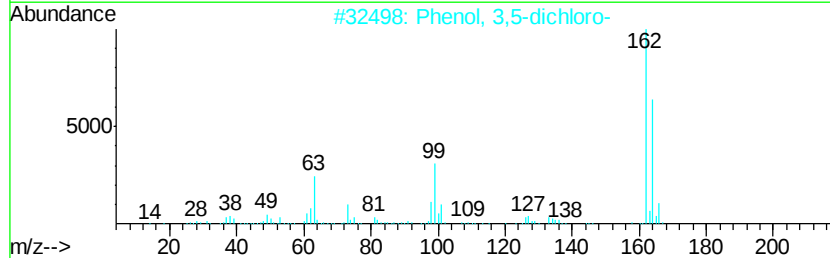
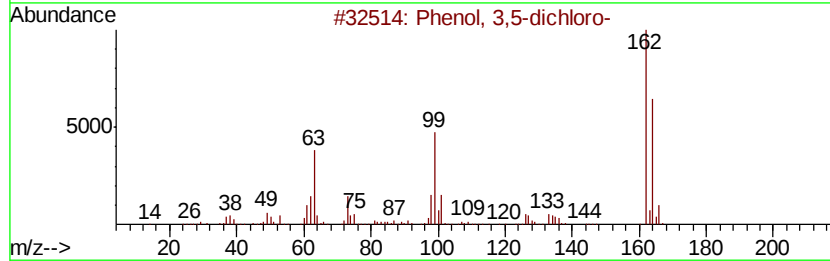
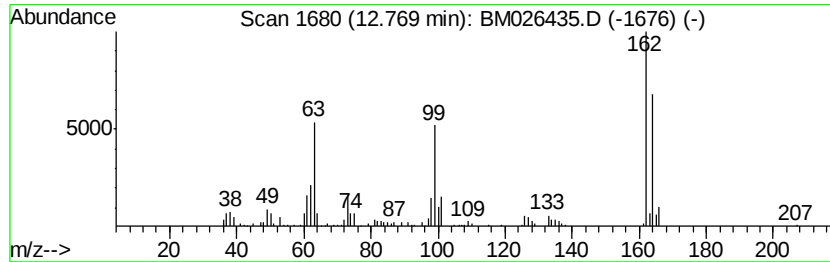
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 Phenol, 3,5-dichloro- Concentration Rank 40

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.77	2.04 ng/ul	197340	Acenaphthene-d10	14.05

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Phenol, 3,5-dichloro-	162	C6H4Cl2O	000591-35-5	98
2		Phenol, 3,5-dichloro-	162	C6H4Cl2O	000591-35-5	97
3		Phenol, 3,4-dichloro-	162	C6H4Cl2O	000095-77-2	97
4		Phenol, 3,4-dichloro-	162	C6H4Cl2O	000095-77-2	94
5		Phenol, 3,5-dichloro-	162	C6H4Cl2O	000591-35-5	94



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM061720\
Data File : BM026435.D
Acq On : 17 Jun 2020 20:43
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Sample : L2959-08
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
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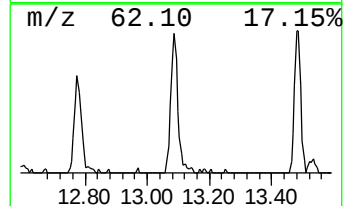
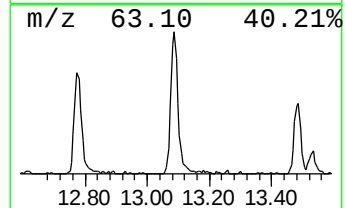
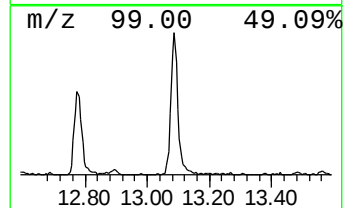
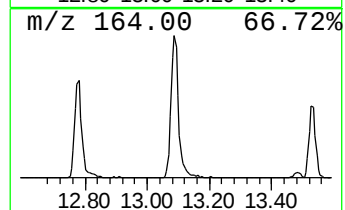
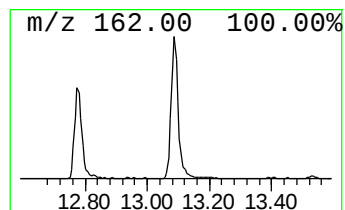
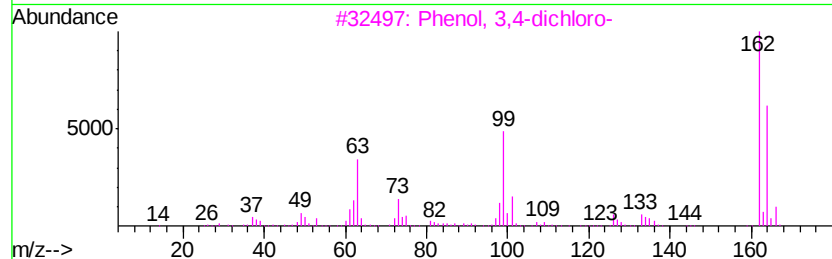
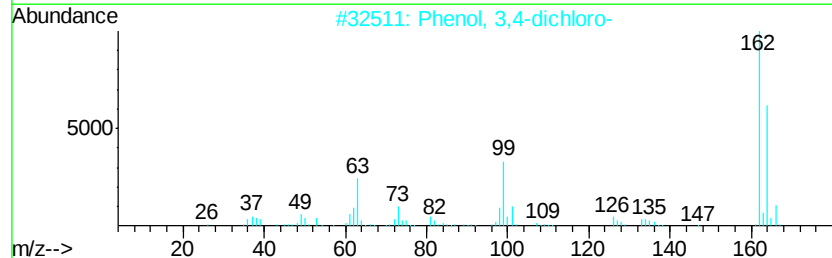
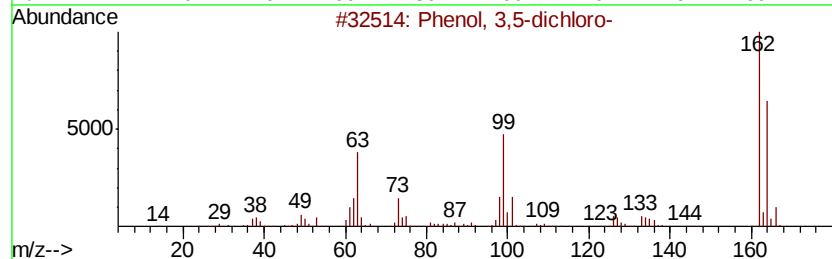
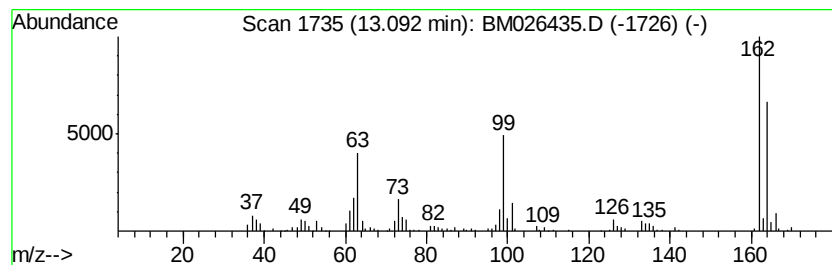
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 Phenol, 3,4-dichloro- Concentration Rank 32

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.09	2.89 ng/ul	279462	Acenaphthene-d10	14.05

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Phenol, 3,5-dichloro-	162	C6H4Cl2O	000591-35-5	98
2		Phenol, 3,4-dichloro-	162	C6H4Cl2O	000095-77-2	97
3		Phenol, 3,4-dichloro-	162	C6H4Cl2O	000095-77-2	97
4		Phenol, 3,4-dichloro-	162	C6H4Cl2O	000095-77-2	96
5		Phenol, 3,5-dichloro-	162	C6H4Cl2O	000591-35-5	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM061720\
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Acq On : 17 Jun 2020 20:43
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Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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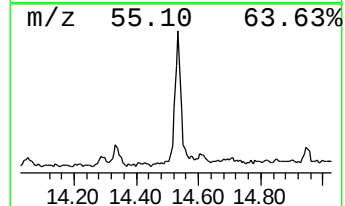
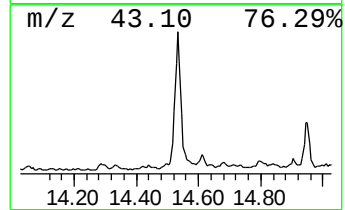
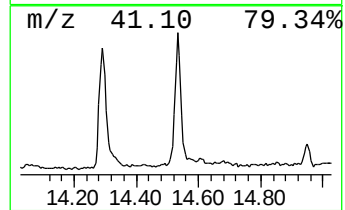
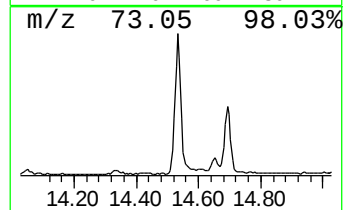
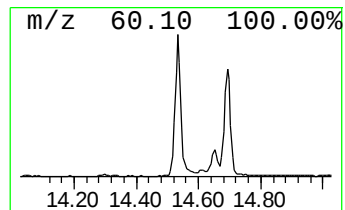
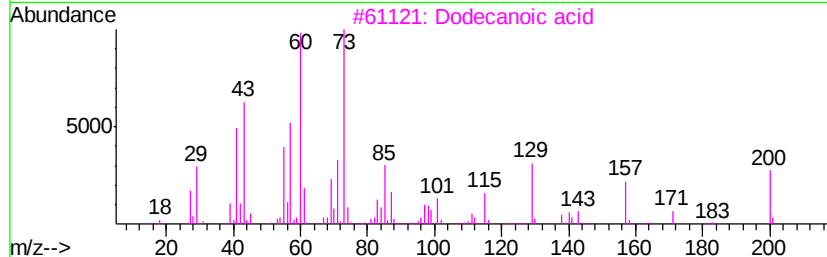
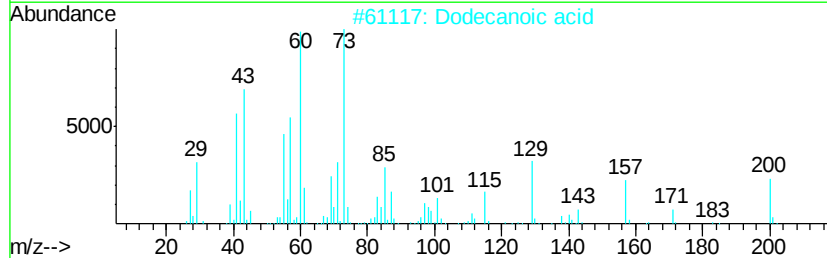
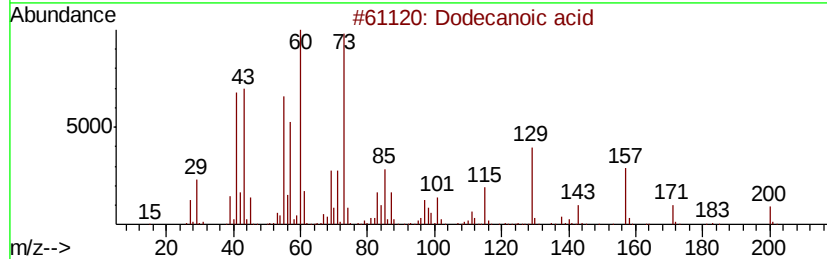
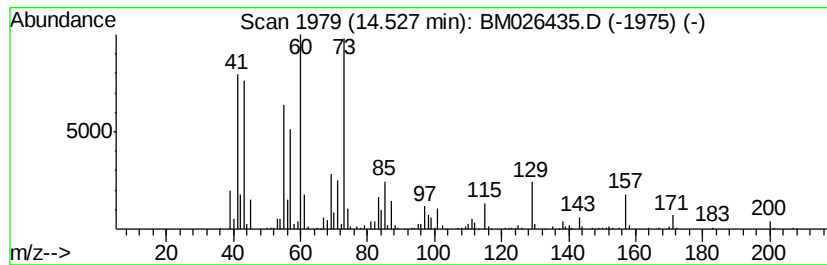
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 Dodecanoic acid Concentration Rank 21

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.53	5.13 ng/ul	495340	Acenaphthene-d10	14.05

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Dodecanoic acid	200	C12H24O2	000143-07-7	97
2			Dodecanoic acid	200	C12H24O2	000143-07-7	91
3			Dodecanoic acid	200	C12H24O2	000143-07-7	91
4			Undecanoic acid	186	C11H22O2	000112-37-8	86
5			Octanoic acid	144	C8H16O2	000124-07-2	83



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM061720\
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Misc :
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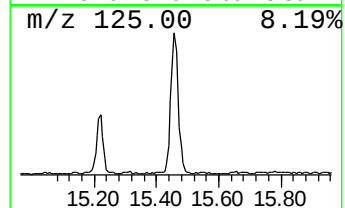
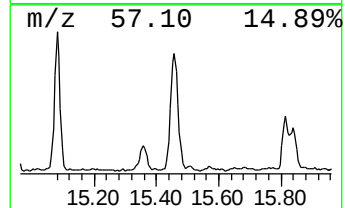
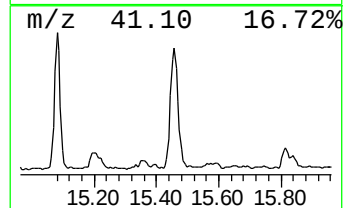
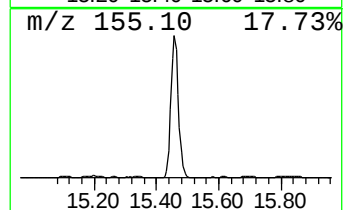
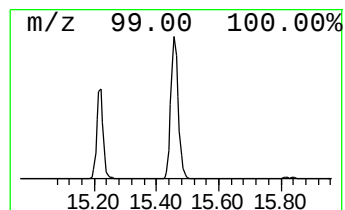
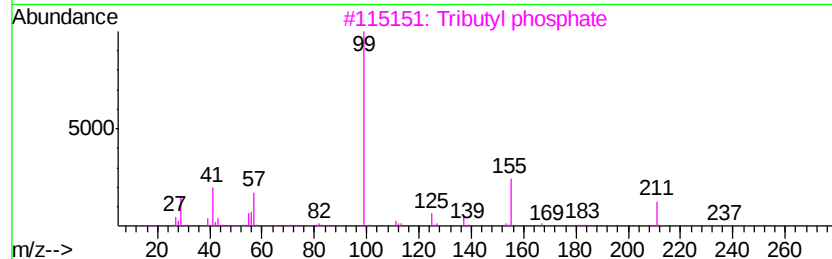
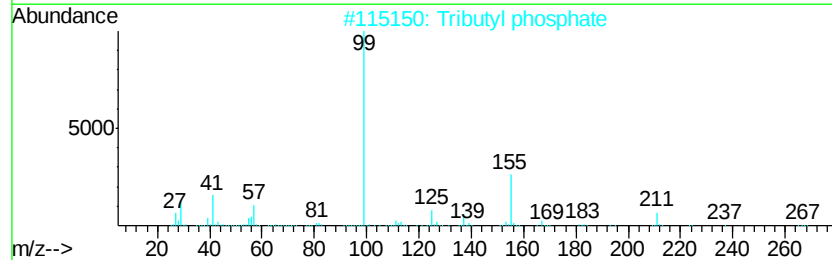
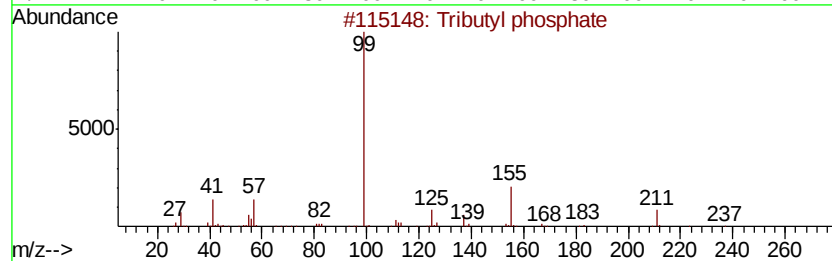
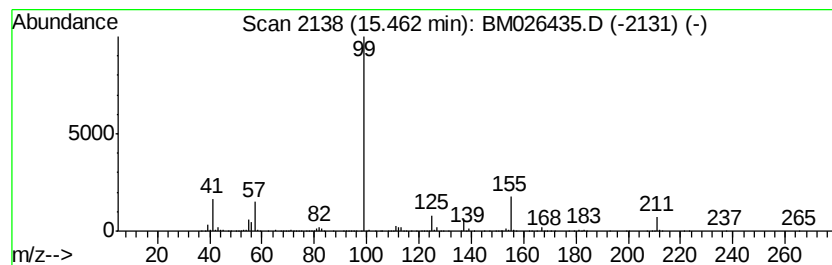
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Quant Title : SVOA CALIBRATION

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

Peak Number 11 Tributyl phosphate Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.46	8.78 ng/ul	1008530	Phenanthrene-d10	16.82

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tributyl phosphate	266	C12H27O4P	000126-73-8	91
2		Tributyl phosphate	266	C12H27O4P	000126-73-8	86
3		Tributyl phosphate	266	C12H27O4P	000126-73-8	83
4		Phosphoric acid, dibutyl 1,1-dim...	352	C13H25F4O4P	1000298-93-9	72
5		Tributyl phosphate	266	C12H27O4P	000126-73-8	58



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Data File : BM026435.D
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Operator : JU/CG
Sample : L2959-08
Misc :
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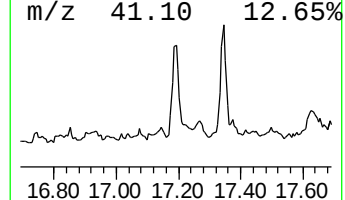
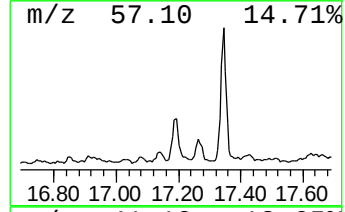
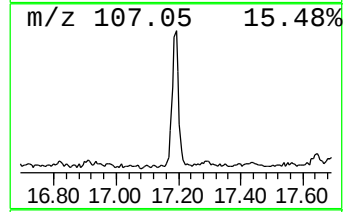
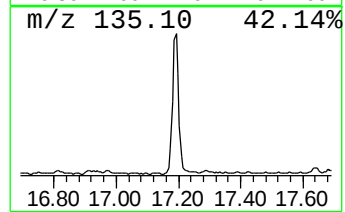
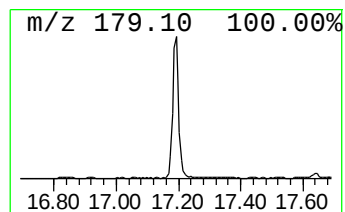
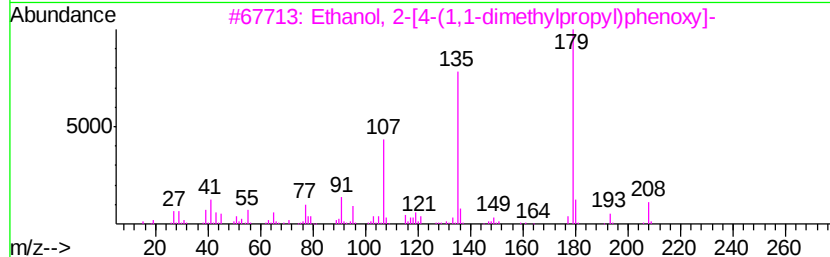
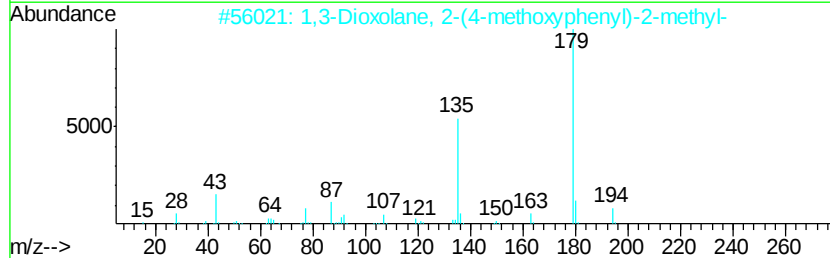
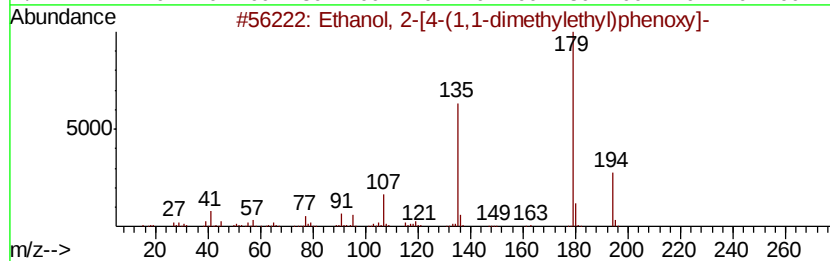
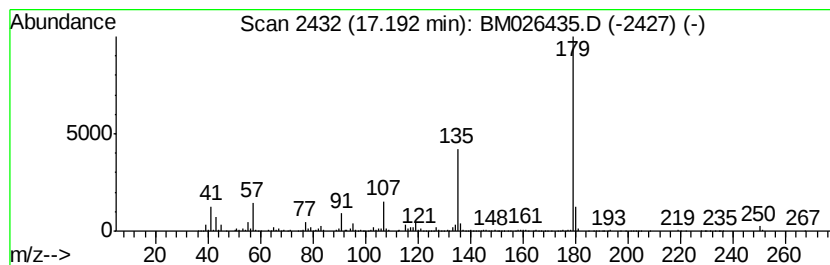
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Quant Title : SVOA CALIBRATION

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

Peak Number 12 Ethanol, 2-[4-(1,1-dimethyl... Concentration Rank 26

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.19	3.93 ng/ul	451018	Phenanthrene-d10	16.82
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Ethanol, 2-[4-(1,1-dimethylethyl)...	194 C12H18O2	000713-46-2 78
2		1,3-Dioxolane, 2-(4-methoxypheny...	194 C11H14O3	036881-00-2 72
3		Ethanol, 2-[4-(1,1-dimethylpropyl)...	208 C13H20O2	006382-07-6 64
4		Benzoic acid, 4-nitroso-, ethyl ...	179 C9H9NO3	007476-79-1 55
5		Acetic acid, [2-[(2-propenylamin...	235 C12H13NO4	000119-45-9 43



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM061720\
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Sample : L2959-08
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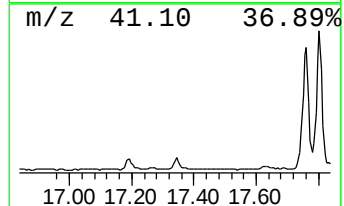
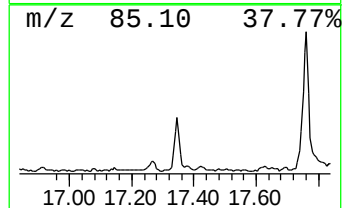
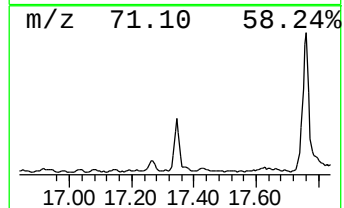
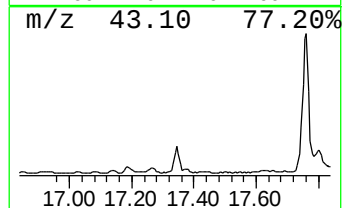
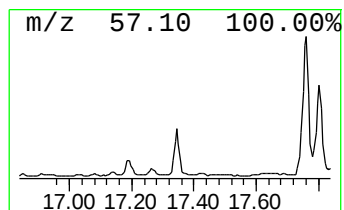
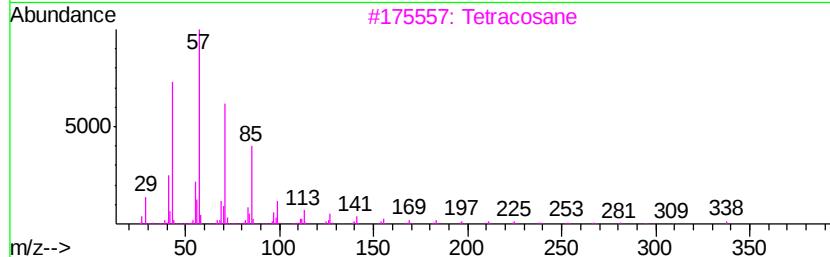
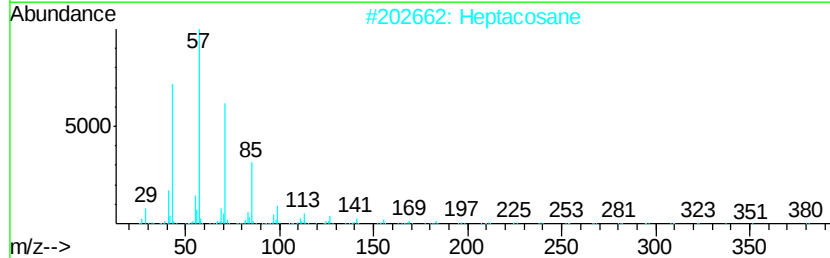
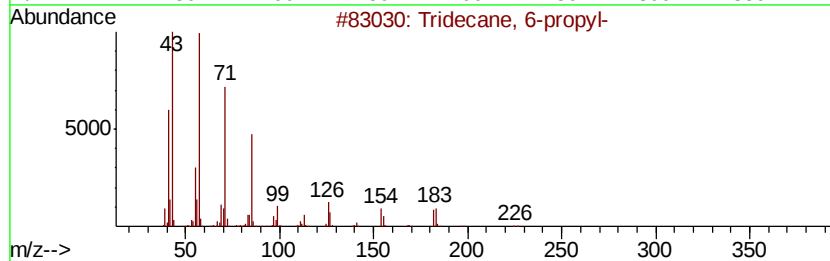
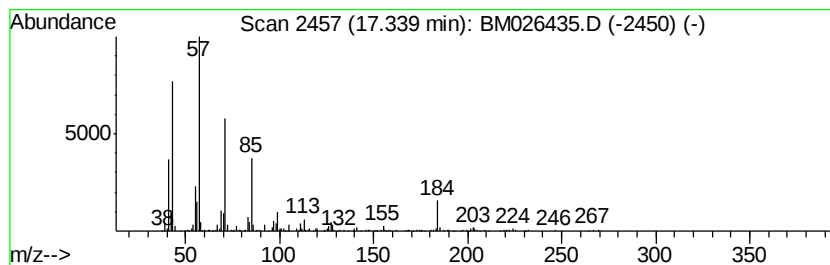
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Quant Title : SVOA CALIBRATION

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

Peak Number 13 (DEL) Alkane: Straight-Chai... Concentration Rank 31

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.34	3.05 ng/ul	350710	Phenanthrene-d10	16.82

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tridecane, 6-propyl-	226	C16H34	055045-10-8	93
2			Heptacosane	380	C27H56	000593-49-7	91
3			Tetracosane	338	C24H50	000646-31-1	91
4			Tridecane, 7-hexyl-	268	C19H40	007225-66-3	91
5			Octacosane	394	C28H58	000630-02-4	91



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Data File : BM026435.D
Acq On : 17 Jun 2020 20:43
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Sample : L2959-08
Misc :
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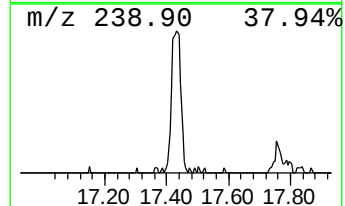
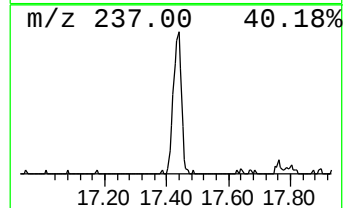
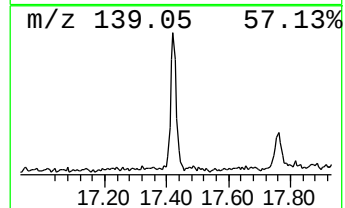
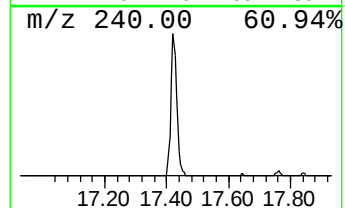
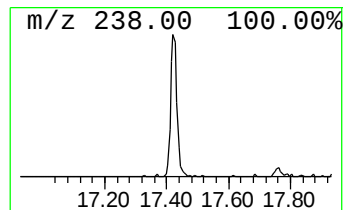
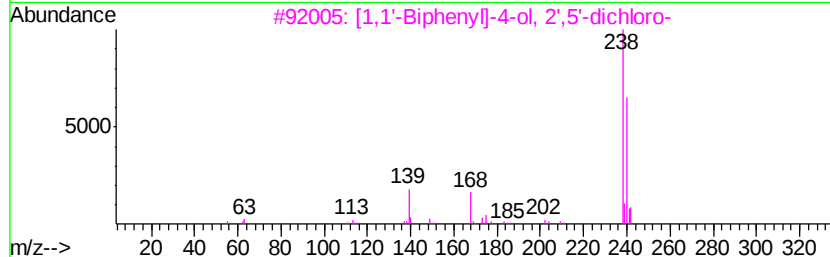
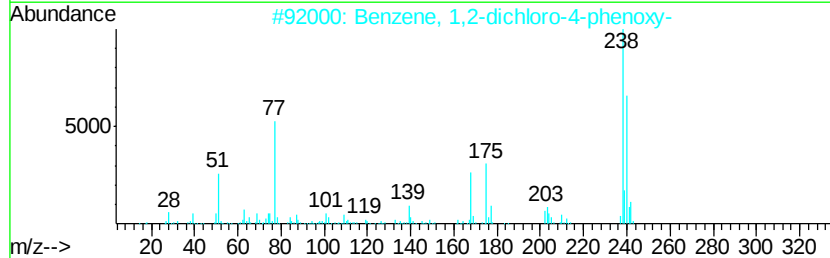
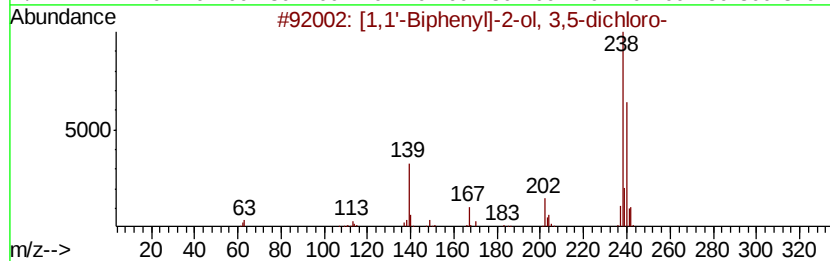
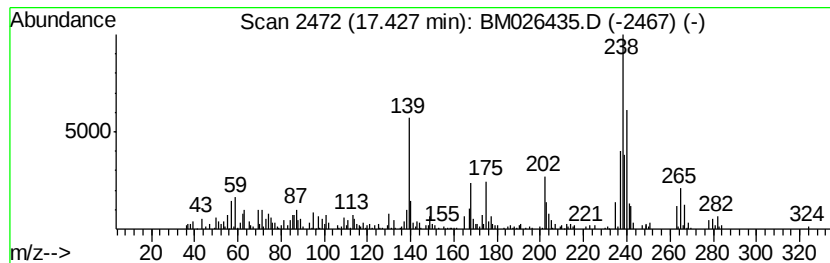
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Quant Title : SVOA CALIBRATION

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

Peak Number 14 [1,1'-Biphenyl]-2-ol, 3,5-d... Concentration Rank 37

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.43	2.21 ng/ul	254264	Phenanthrene-d10	16.82

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			[1,1'-Biphenyl]-2-ol, 3,5-dichloro-	238	C12H8Cl2O	005335-24-0	86
2			Benzene, 1,2-dichloro-4-phenoxy-	238	C12H8Cl2O	055538-69-7	70
3			[1,1'-Biphenyl]-4-ol, 2',5'-dich...	238	C12H8Cl2O	053905-28-5	60
4			[1,1'-Biphenyl]-4-ol, 3,5-dichloro-	238	C12H8Cl2O	001137-59-3	55
5			[1,1'-Biphenyl]-4-ol, 3,4'-dichl...	238	C12H8Cl2O	053905-31-0	55



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM061720\
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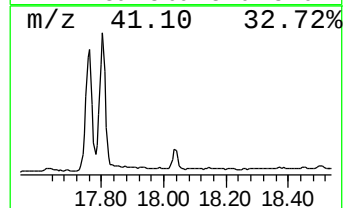
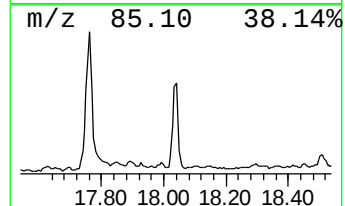
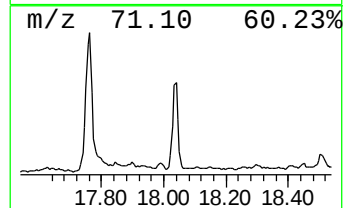
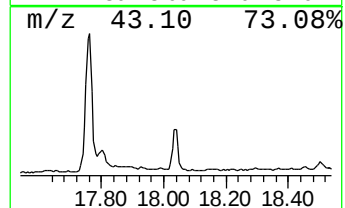
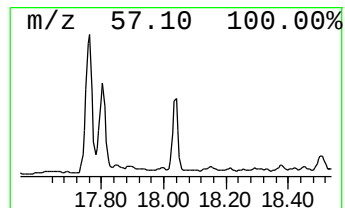
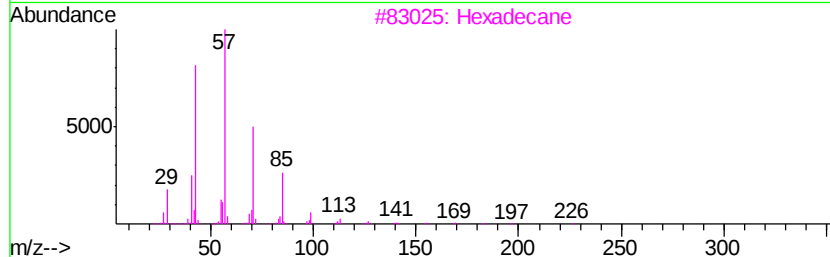
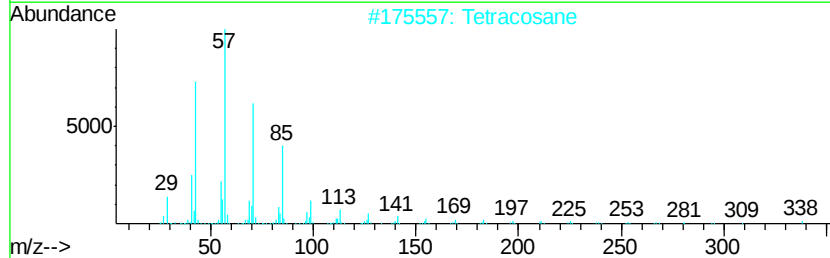
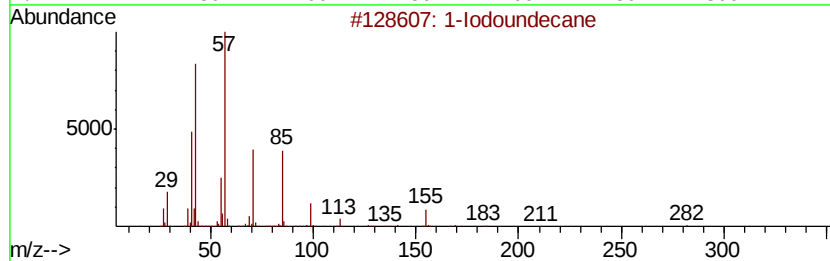
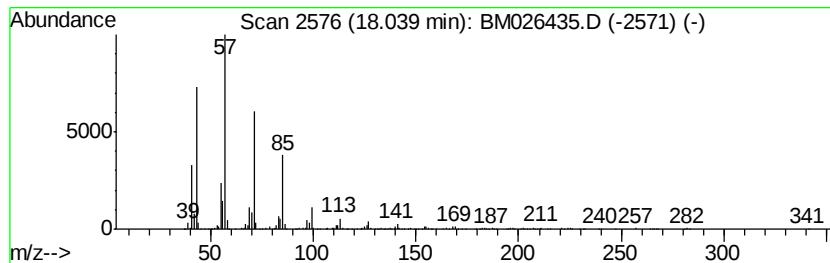
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Quant Title : SVOA CALIBRATION

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

Peak Number 15 1-Iodoundecane Concentration Rank 27

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.04	3.45 ng/ul	396728	Phenanthrene-d10	16.82

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Iodoundecane	282	C11H23I	004282-44-4	92
2			Tetracosane	338	C24H50	000646-31-1	91
3			Hexadecane	226	C16H34	000544-76-3	90
4			Dodecane, 2-methyl-6-propyl-	226	C16H34	055045-08-4	90
5			Tridecane, 1-iodo-	310	C13H27I	035599-77-0	90



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM061720\
Data File : BM026435.D
Acq On : 17 Jun 2020 20:43
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Misc :
ALS Vial : 11 Sample Multiplier: 1

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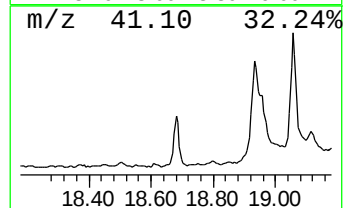
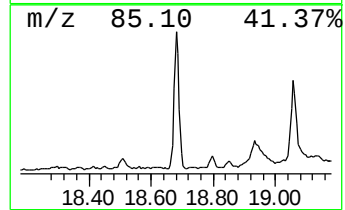
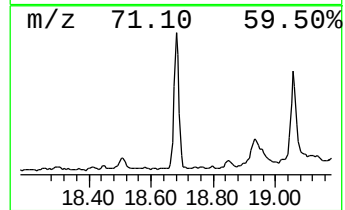
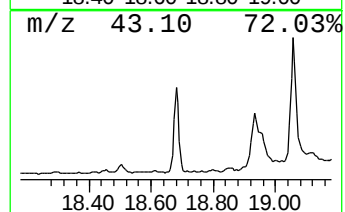
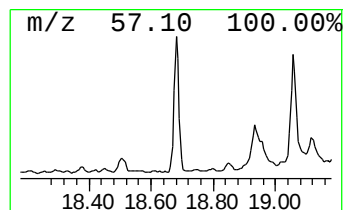
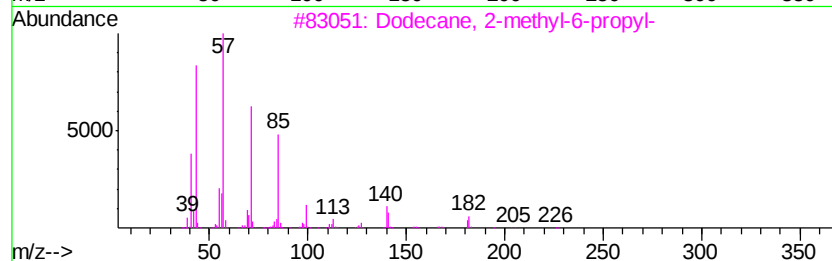
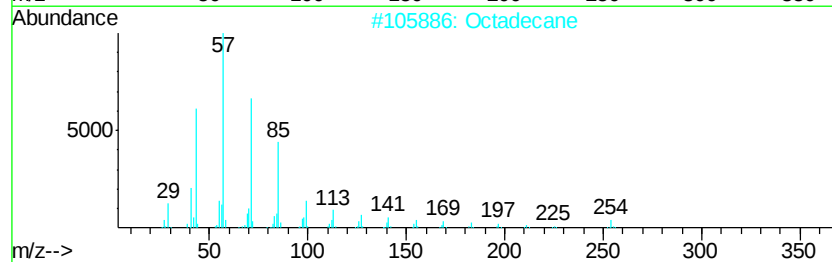
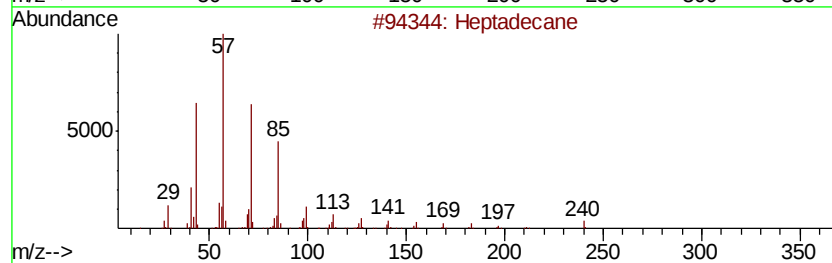
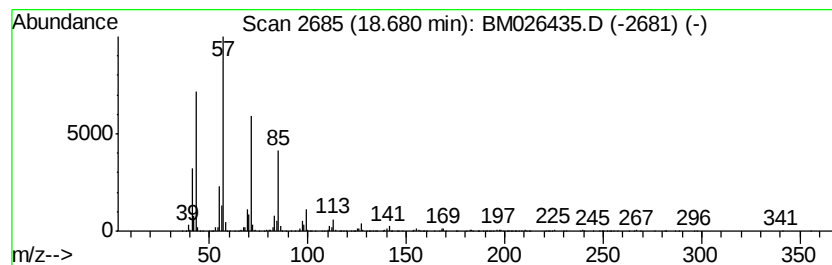
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Quant Title : SVOA CALIBRATION

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

Peak Number 16 (DEL) Alkane: Straight-Chai... Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.68	6.57 ng/ul	754187	Phenanthrene-d10	16.82

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heptadecane	240	C17H36	000629-78-7	95
2			Octadecane	254	C18H38	000593-45-3	95
3			Dodecane, 2-methyl-6-propyl-	226	C16H34	055045-08-4	90
4			Eicosane, 10-methyl-	296	C21H44	054833-23-7	90
5			Nonadecane	268	C19H40	000629-92-5	90



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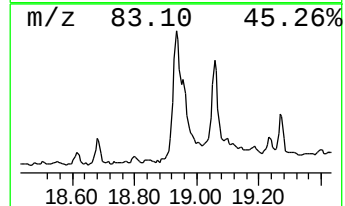
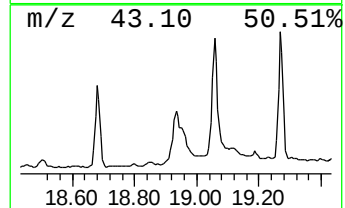
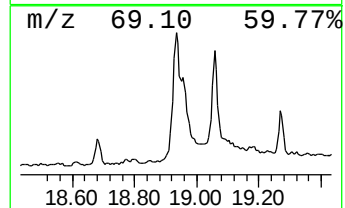
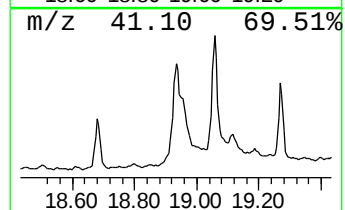
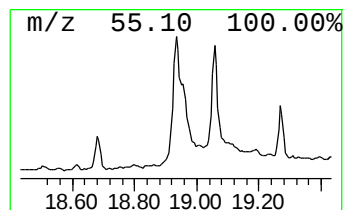
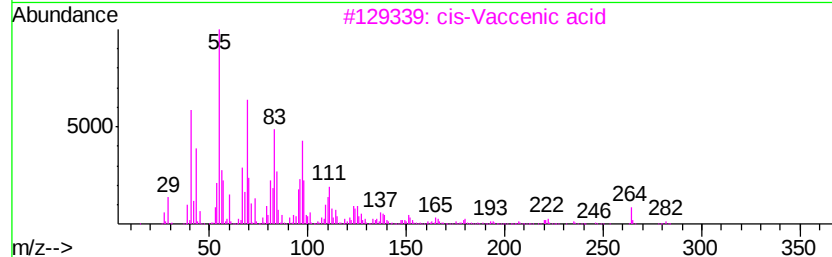
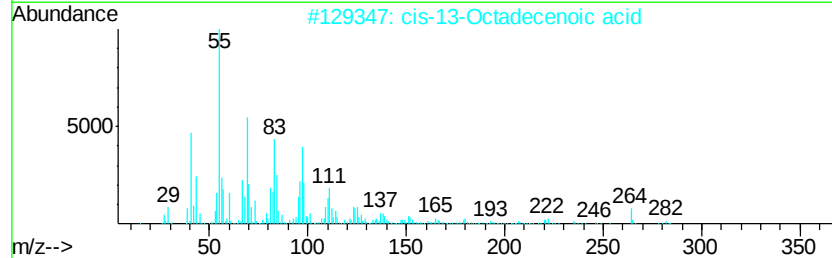
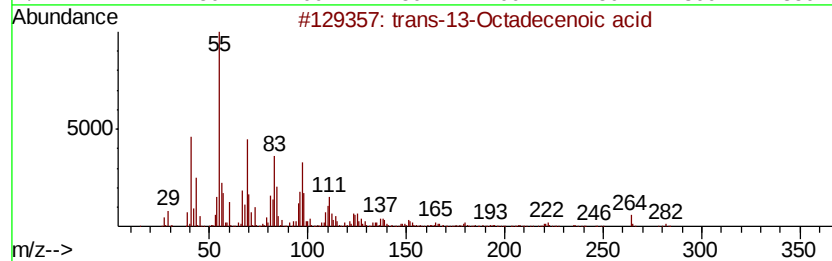
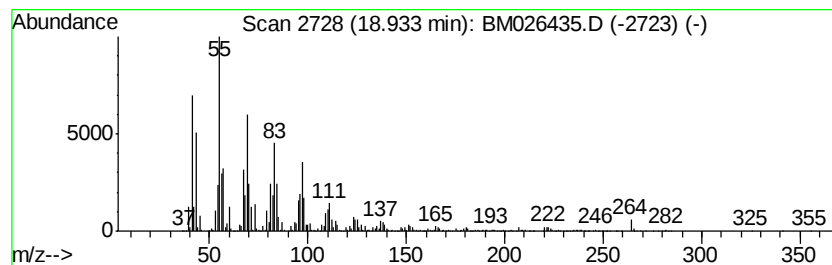
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Quant Title : SVOA CALIBRATION

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

Peak Number 17 trans-13-Octadecenoic acid Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.93	22.09 ng/ul	2983400	Chrysene-d12	21.02

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	trans-13-Octadecenoic acid	282	C18H34O2	000693-71-0	98
2		cis-13-Octadecenoic acid	282	C18H34O2	013126-39-1	96
3		cis-Vaccenic acid	282	C18H34O2	000506-17-2	95
4		6-Octadecenoic acid	282	C18H34O2	1000336-66-8	95
5		Oleic Acid	282	C18H34O2	000112-80-1	94



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Data File : BM026435.D
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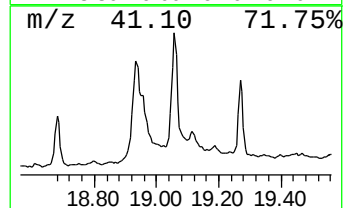
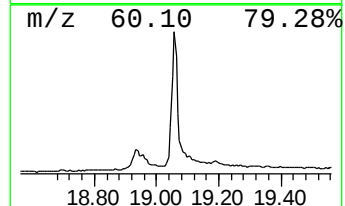
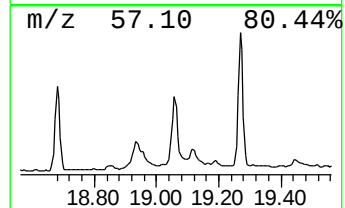
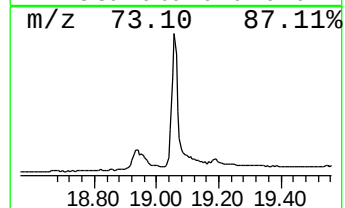
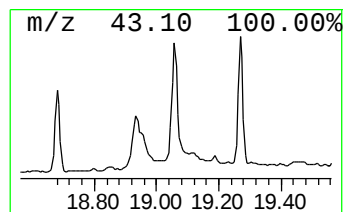
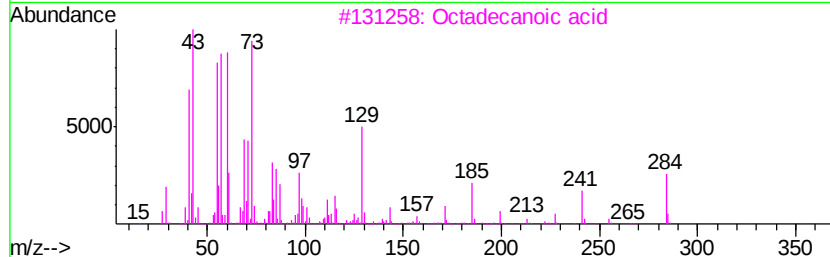
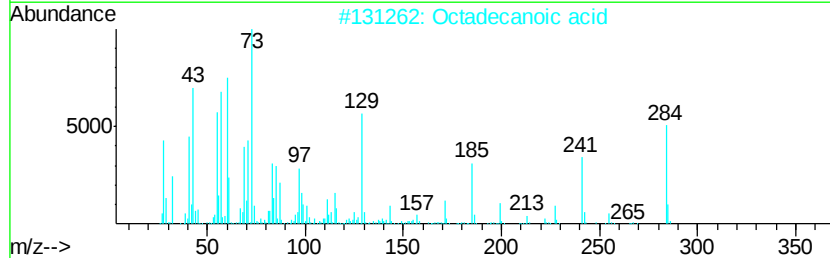
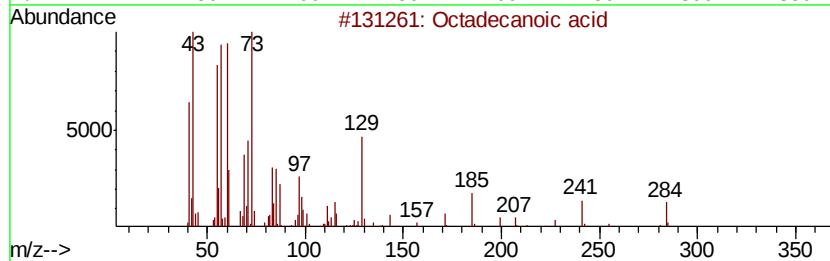
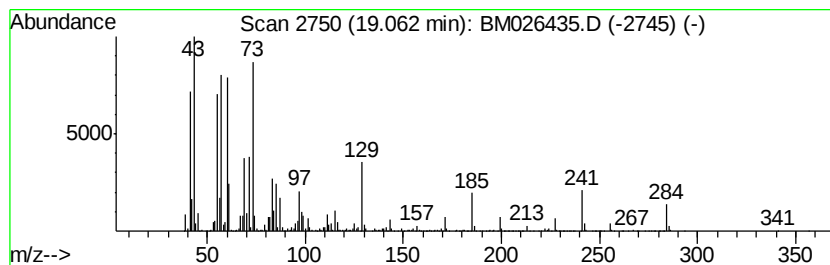
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Quant Title : SVOA CALIBRATION

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

Peak Number 18 Octadecanoic acid Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.06	14.48 ng/ul	1956050	Chrysene-d12	21.02

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octadecanoic acid	284	C18H36O2	000057-11-4	99
2		Octadecanoic acid	284	C18H36O2	000057-11-4	99
3		Octadecanoic acid	284	C18H36O2	000057-11-4	99
4		Octadecanoic acid	284	C18H36O2	000057-11-4	99
5		Octadecanoic acid	284	C18H36O2	000057-11-4	93



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM061720\
Data File : BM026435.D
Acq On : 17 Jun 2020 20:43
Operator : JU/CG
Sample : L2959-08
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
BG2K2

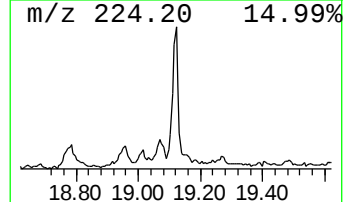
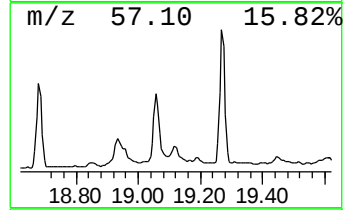
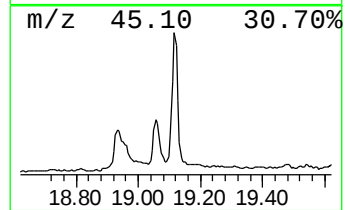
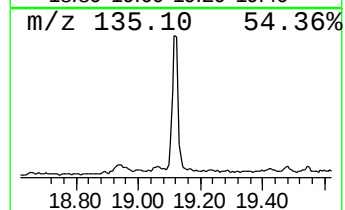
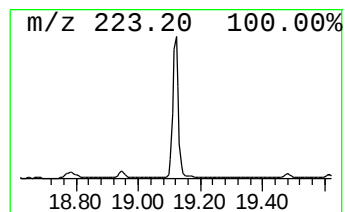
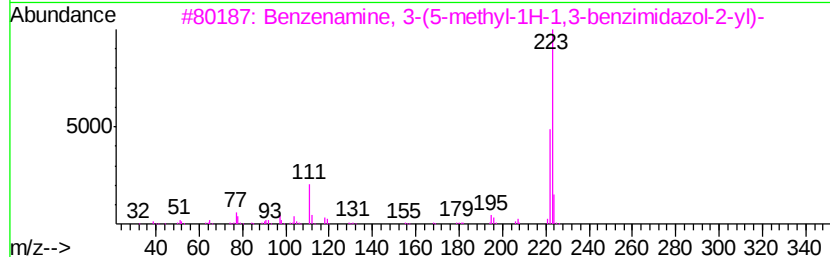
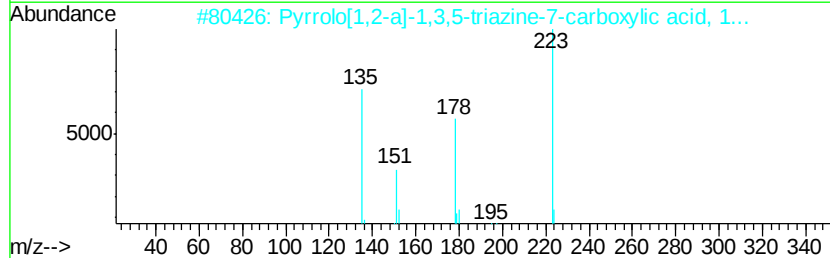
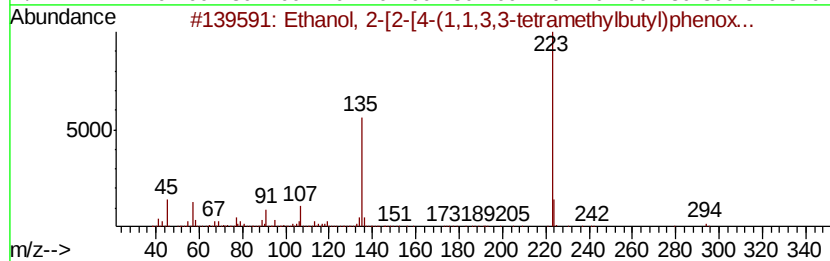
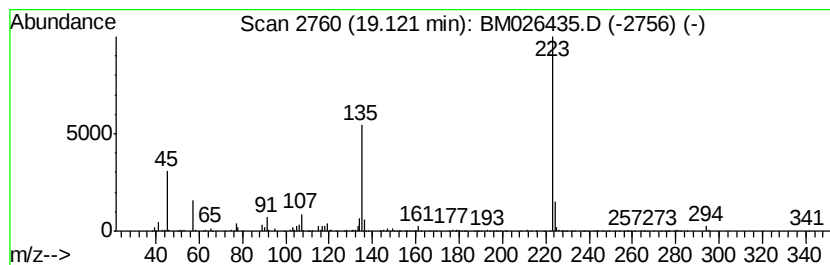
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Quant Title : SVOA CALIBRATION

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

Peak Number 19 Ethanol, 2-[2-[4-(1,1,3,3-t... Concentration Rank 23

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.12	4.72 ng/ul	637047	Chrysene-d12	21.02

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Ethanol, 2-[2-[4-(1,1,3,3-tetram...	294	C18H30O3	002315-61-9	95
2		Pyrrolo[1,2-a]-1,3,5-triazine-7-...	223	C9H9N3O4	054449-89-7	90
3		Benzenamine, 3-(5-methyl-1H-1,3-...	223	C14H13N3	1000319-43-9	43
4		Tricyclo[3.3.3.0(1,5)]undec-6-en...	223	C14H13N3	118894-71-6	43
5		Tetrazole, 1-(4-ethylphenyl)-5-(...	410	C25H26N6	125038-06-4	32



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM061720\
Data File : BM026435.D
Acq On : 17 Jun 2020 20:43
Operator : JU/CG
Sample : L2959-08
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
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ClientSampleId :
BG2K2

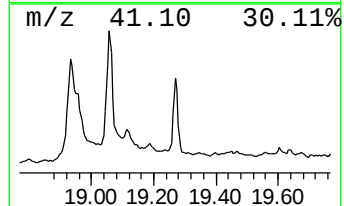
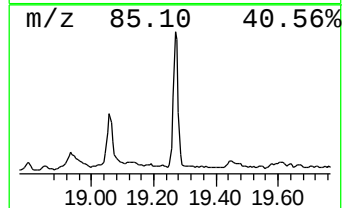
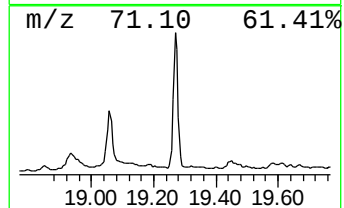
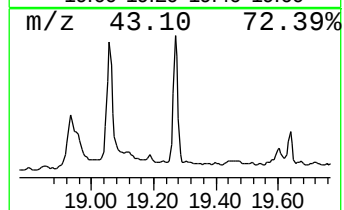
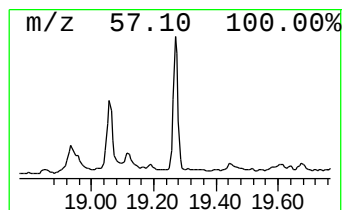
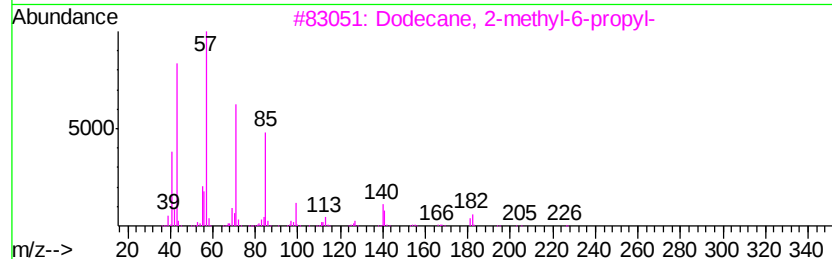
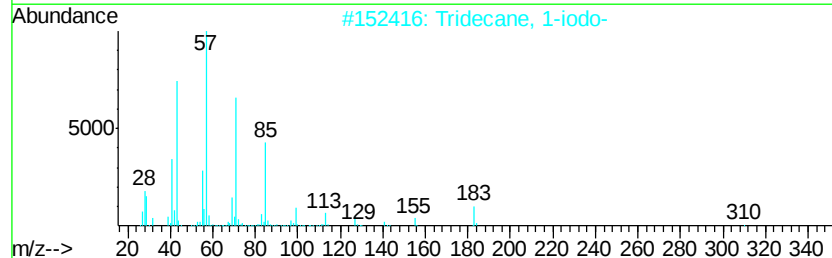
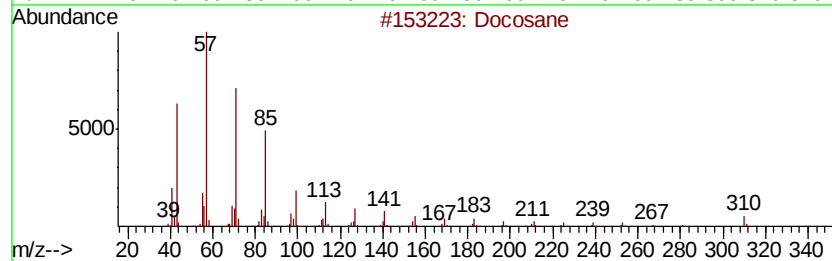
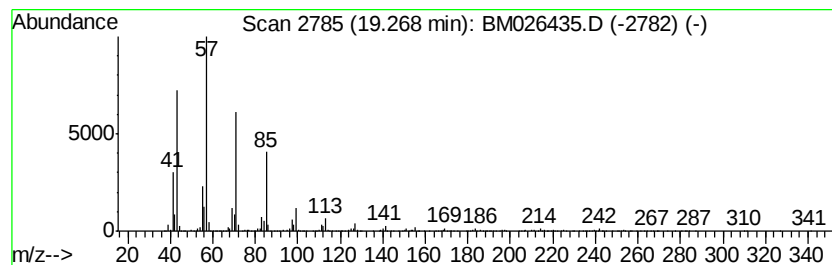
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Quant Title : SVOA CALIBRATION

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

Peak Number 20 (DEL) Alkane: Straight-Chai... Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.27	8.55 ng/ul	1154830	Chrysene-d12	21.02

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Docosane	310	C22H46	000629-97-0	97
2			Tridecane, 1-iodo-	310	C13H27I	035599-77-0	94
3			Dodecane, 2-methyl-6-propyl-	226	C16H34	055045-08-4	93
4			Heptacosane	380	C27H56	000593-49-7	91
5			Heneicosane	296	C21H44	000629-94-7	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM061720\
Data File : BM026435.D
Acq On : 17 Jun 2020 20:43
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Sample : L2959-08
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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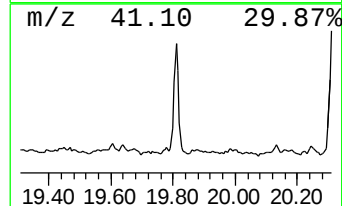
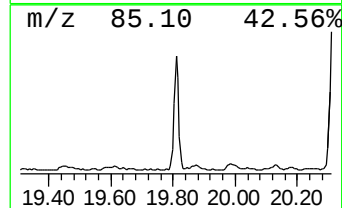
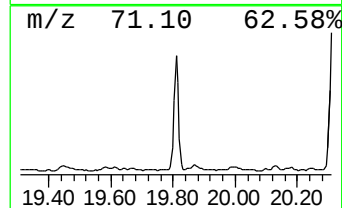
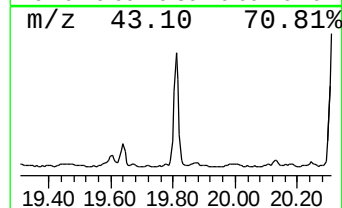
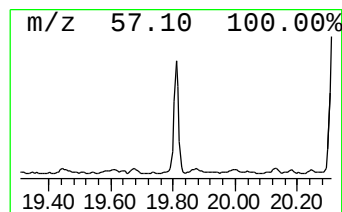
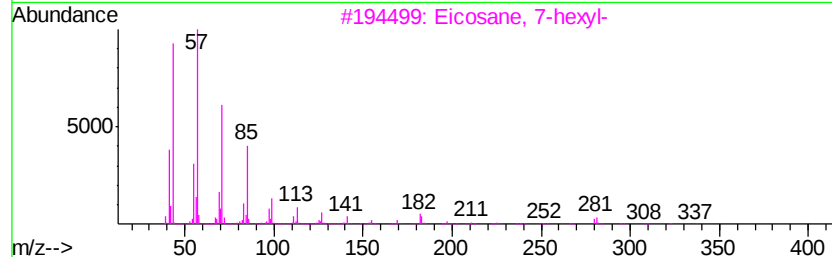
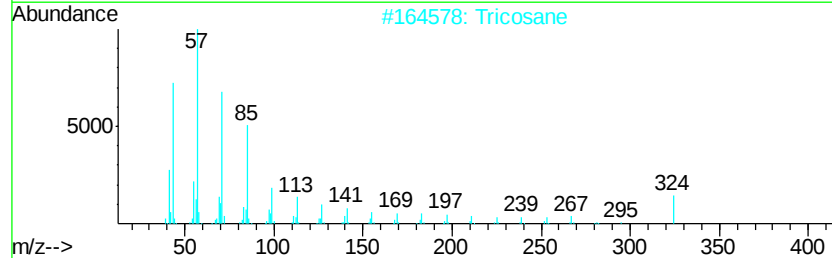
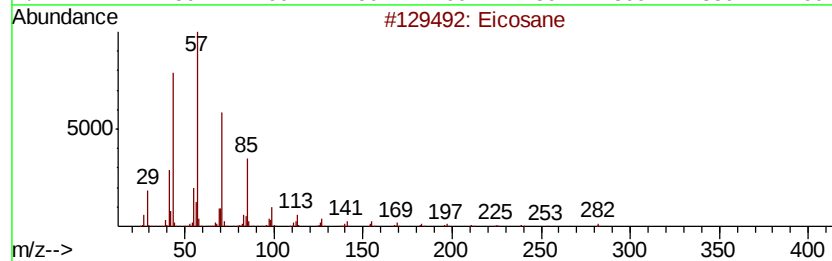
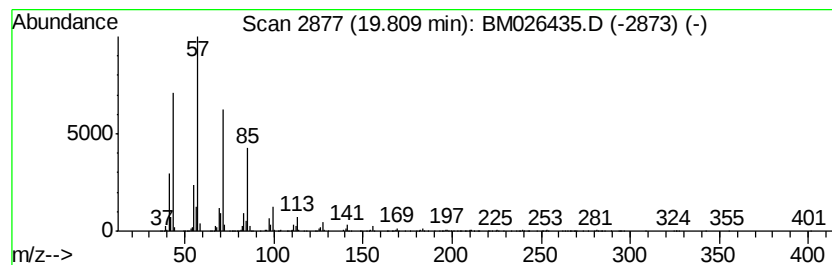
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Quant Title : SVOA CALIBRATION

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

Peak Number 21 (DEL) Alkane: Straight-Chai... Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.81	10.40 ng/ul	1404850	Chrysene-d12	21.02

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Eicosane	282	C20H42	000112-95-8	95
2		Tricosane	324	C23H48	000638-67-5	94
3		Eicosane, 7-hexyl-	366	C26H54	055333-99-8	91
4		Dodecane, 2-methyl-6-propyl-	226	C16H34	055045-08-4	91
5		Heptacosane	380	C27H56	000593-49-7	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM061720\
Data File : BM026435.D
Acq On : 17 Jun 2020 20:43
Operator : JU/CG
Sample : L2959-08
Misc :
ALS Vial : 11 Sample Multiplier: 1

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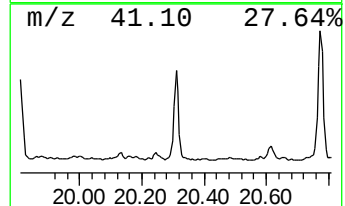
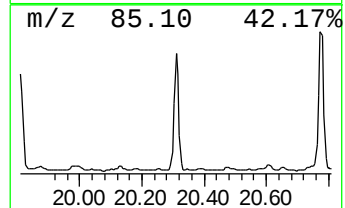
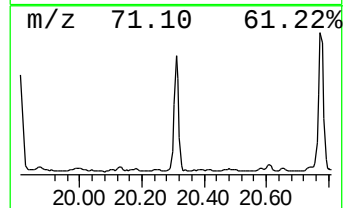
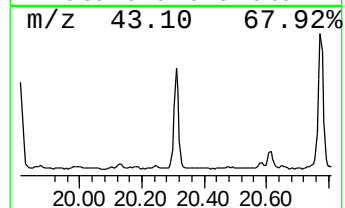
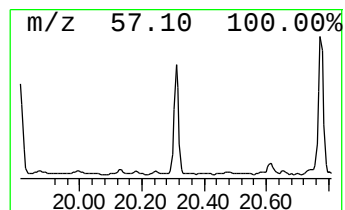
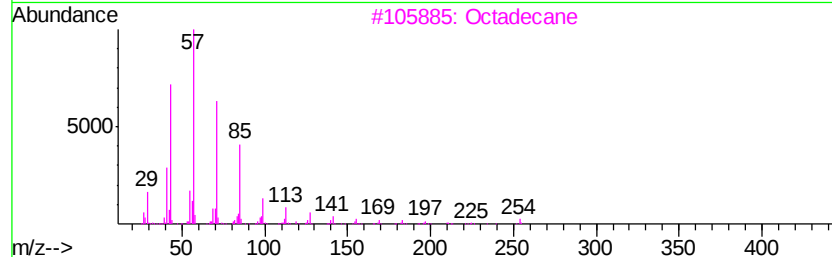
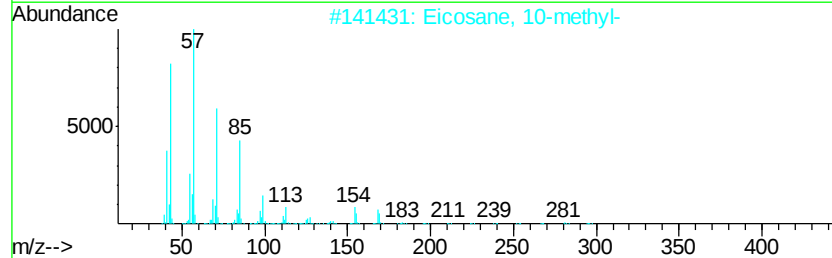
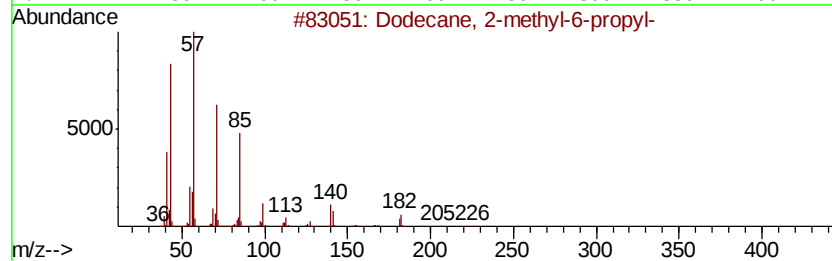
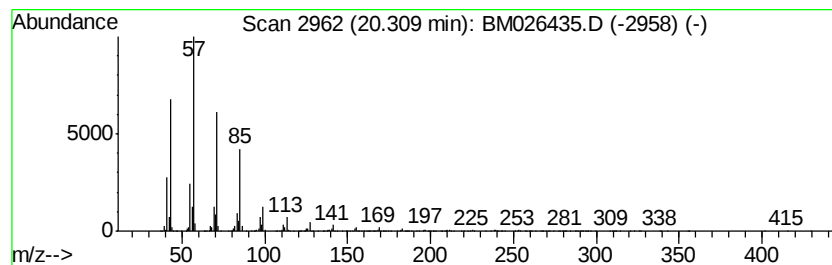
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Quant Title : SVOA CALIBRATION

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

Peak Number 22 (DEL) Alkane: Straight-Chai... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.31	12.74 ng/ul	1721090	Chrysene-d12	21.02

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Dodecane, 2-methyl-6-propyl-	226	C16H34	055045-08-4	94
2		Eicosane, 10-methyl-	296	C21H44	054833-23-7	94
3		Octadecane	254	C18H38	000593-45-3	93
4		Heptacosane	380	C27H56	000593-49-7	91
5		Eicosane	282	C20H42	000112-95-8	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM061720\
Data File : BM026435.D
Acq On : 17 Jun 2020 20:43
Operator : JU/CG
Sample : L2959-08
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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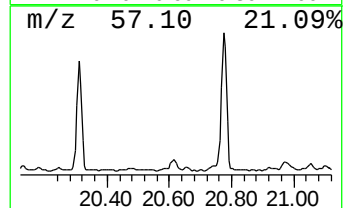
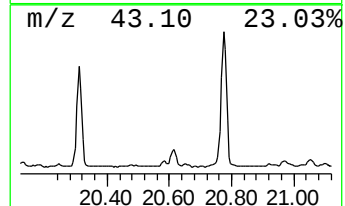
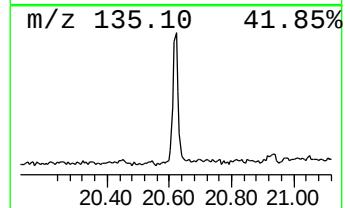
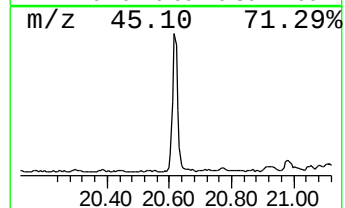
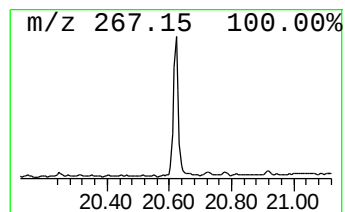
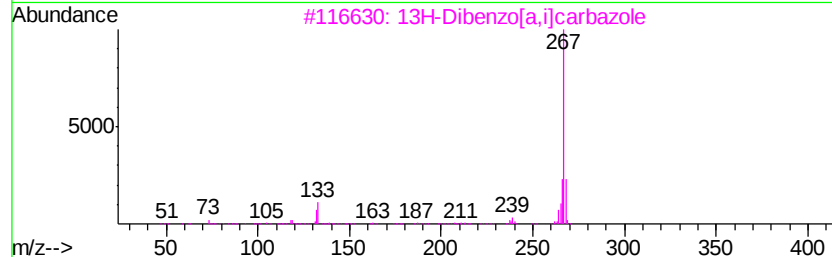
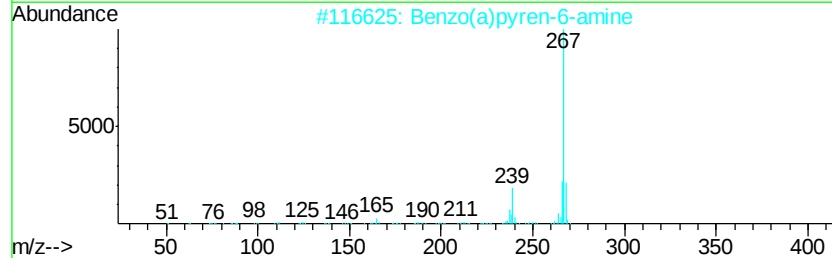
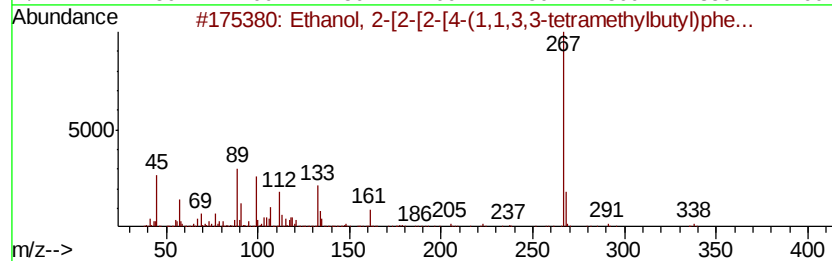
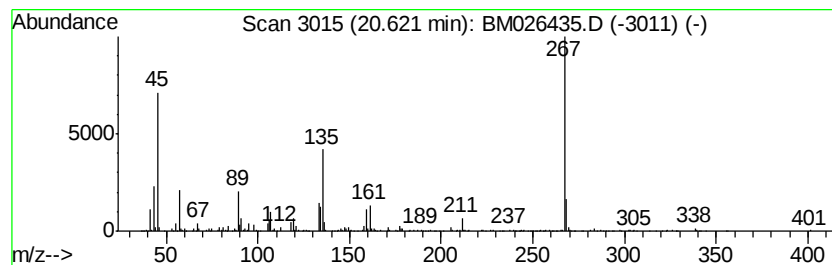
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Quant Title : SVOA CALIBRATION

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

Peak Number 23 Ethanol, 2-[2-[2-[4-(1,1,3,... Concentration Rank 25

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.62	4.14 ng/ul	559610	Chrysene-d12	21.02

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Ethanol, 2-[2-[2-[4-(1,1,3,3-tet...	338	C20H34O4	002315-62-0	70
2		Benzo(a)pyren-6-amine	267	C20H13N	007428-83-3	27
3		13H-Dibenzo[a,i]carbazole	267	C20H13N	000239-64-5	25
4		7H-Dibenzo(a,q)carbazole	267	C20H13N	000207-84-1	25
5		13H-Dibenzo[a,i]carbazole	267	C20H13N	000239-64-5	25



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM061720\
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Acq On : 17 Jun 2020 20:43
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Sample : L2959-08
Misc :
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Instrument :
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ClientSampled :
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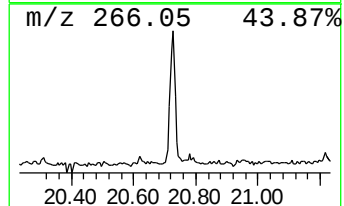
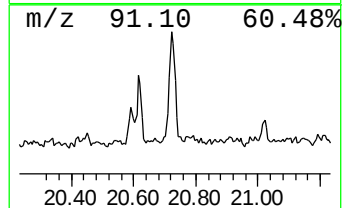
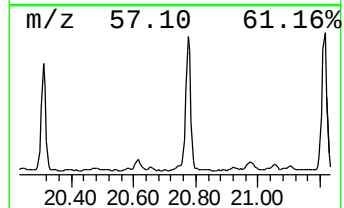
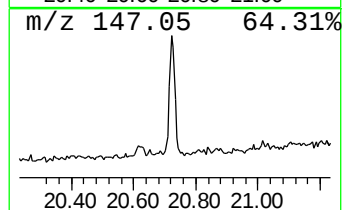
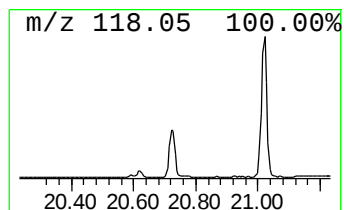
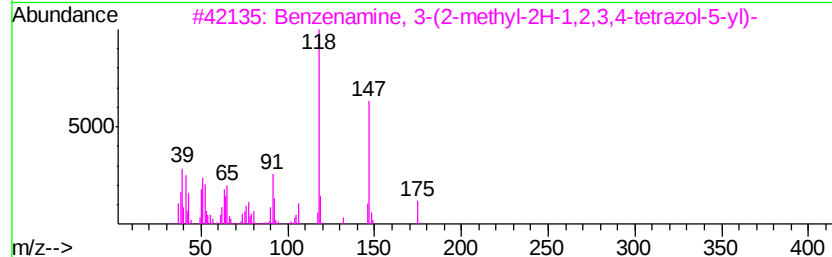
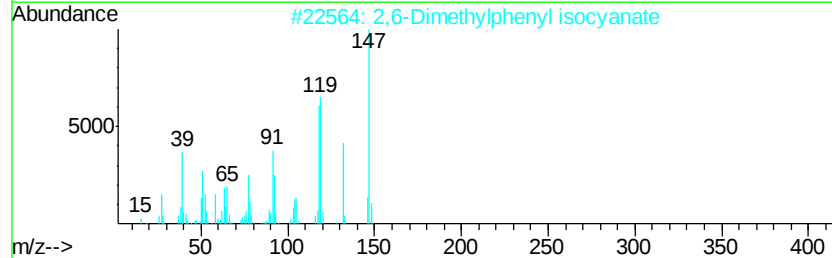
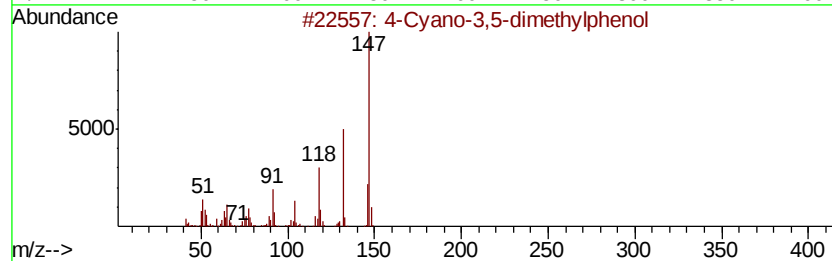
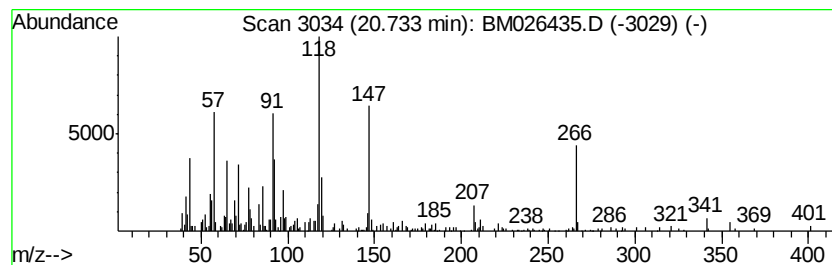
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Quant Title : SVOA CALIBRATION

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

Peak Number 24 unknown-01 Concentration Rank 39

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.73	2.12 ng/ul	286364	Chrysene-d12	21.02

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	4-Cyano-3,5-dimethylphenol	147	C9H9NO	058537-99-8	40
2		2,6-Dimethylphenyl isocyanate	147	C9H9NO	028556-81-2	18
3		Benzenamine, 3-(2-methyl-2H-1,2,4-triazol-5-yl)-	175	C8H9N5	1000362-62-4	16
4		1H-Indole-1-carboxaldehyde, 2,3-dimethyl-	147	C9H9NO	002861-59-8	16
5		Benzenepropanamine	135	C9H13N	002038-57-5	14



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM061720\
Data File : BM026435.D
Acq On : 17 Jun 2020 20:43
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Sample : L2959-08
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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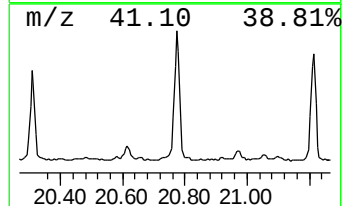
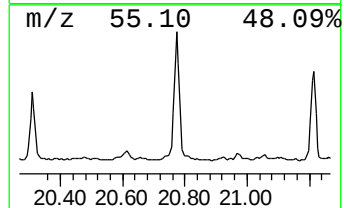
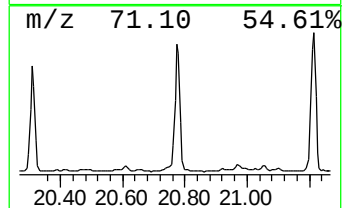
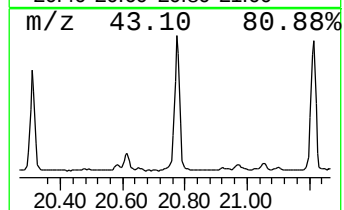
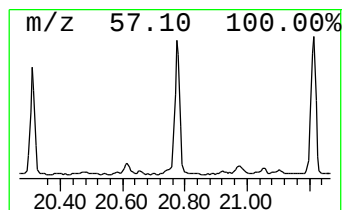
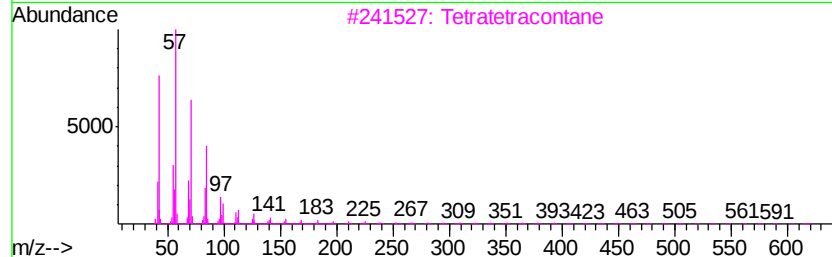
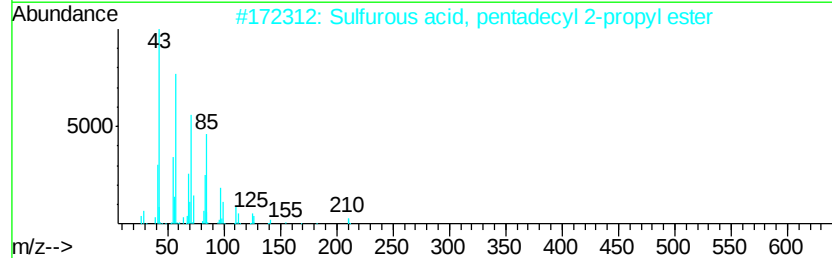
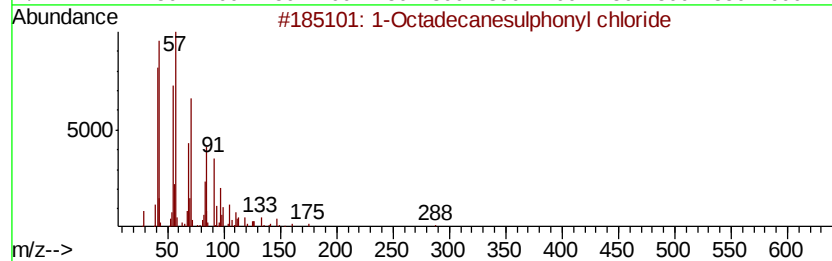
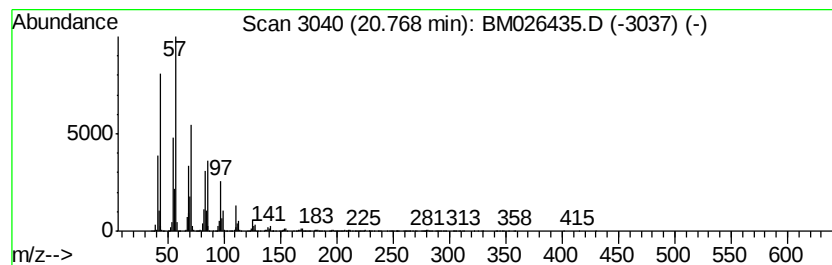
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Quant Title : SVOA CALIBRATION

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

Peak Number 25 1-Octadecanesulphonyl chloride Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.77	20.78 ng/ul	2806570	Chrysene-d12	21.02

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1-Octadecanesulphonyl chloride	352	C18H37ClO2S	1000342-70-4	91
2		Sulfurous acid, pentadecyl 2-pro...	334	C18H38O3S	1000309-12-6	91
3		Tetratetracontane	619	C44H90	007098-22-8	87
4		10-Methylnonadecane	282	C20H42	056862-62-5	83
5		Sulfurous acid, butyl heptadecyl...	376	C21H44O3S	1000309-18-4	83



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM061720\
Data File : BM026435.D
Acq On : 17 Jun 2020 20:43
Operator : JU/CG
Sample : L2959-08
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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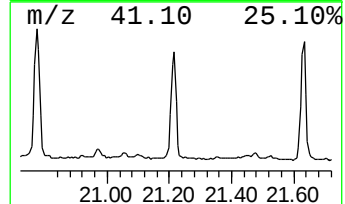
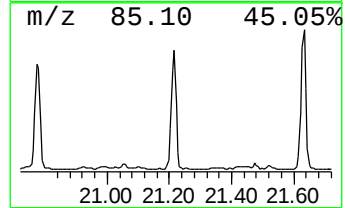
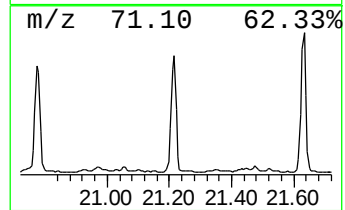
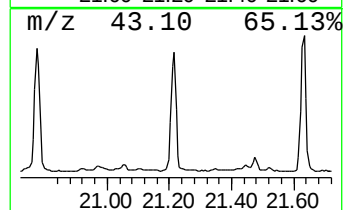
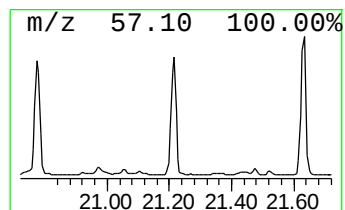
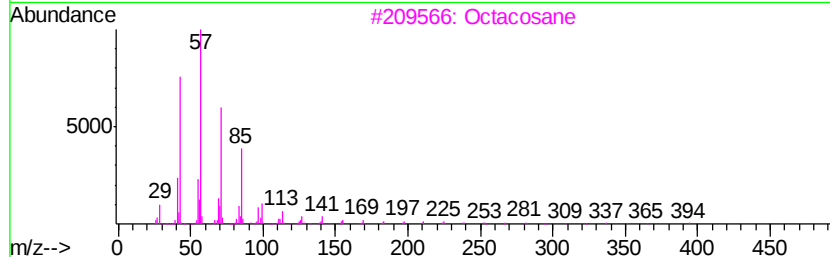
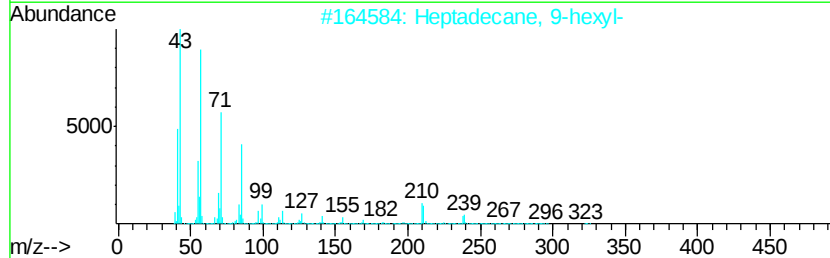
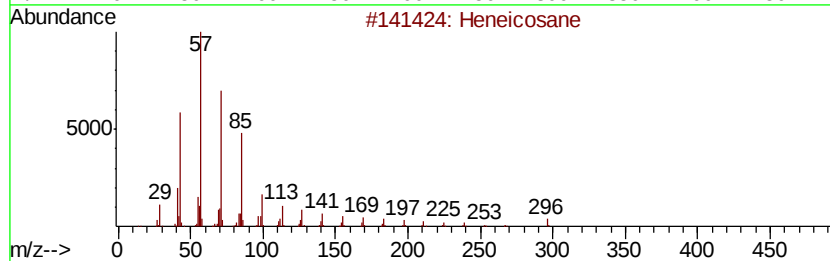
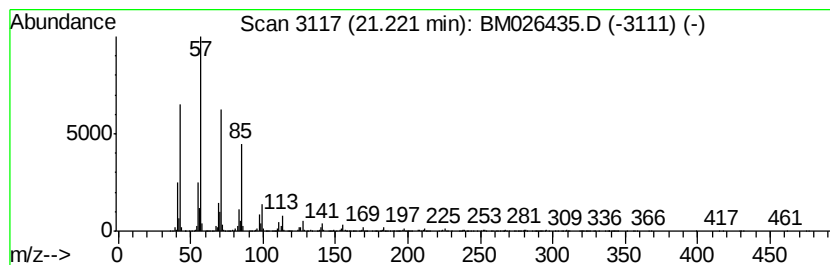
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Quant Title : SVOA CALIBRATION

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

Peak Number 26 (DEL) Alkane: Straight-Chai... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.22	16.95 ng/ul	2289650	Chrysene-d12	21.02

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heneicosane	296	C21H44	000629-94-7	96
2		Heptadecane, 9-hexyl-	324	C23H48	055124-79-3	94
3		Octacosane	394	C28H58	000630-02-4	91
4		Heneicosane	296	C21H44	000629-94-7	91
5		Octacosane	394	C28H58	000630-02-4	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM061720\
Data File : BM026435.D
Acq On : 17 Jun 2020 20:43
Operator : JU/CG
Sample : L2959-08
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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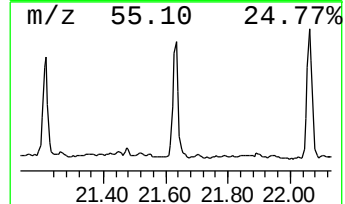
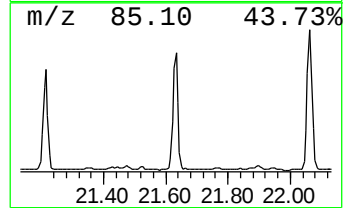
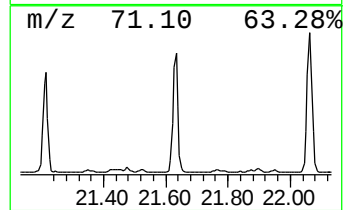
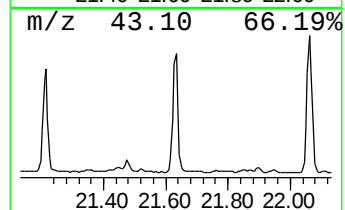
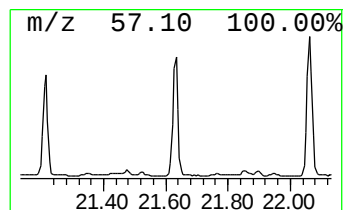
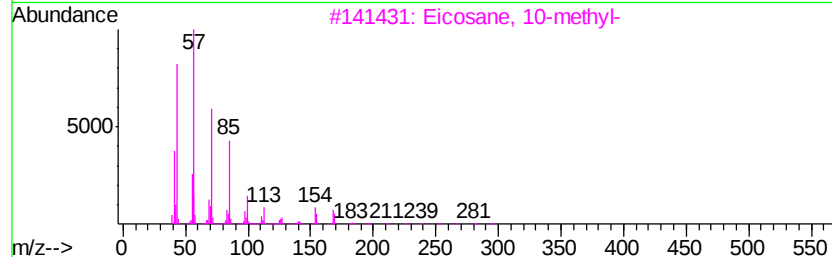
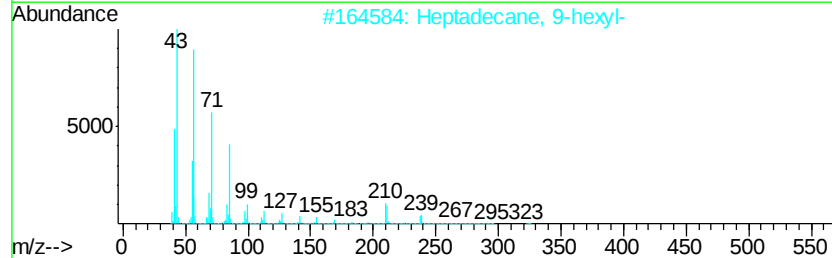
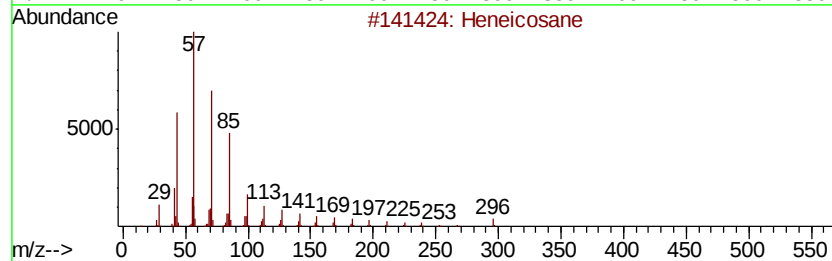
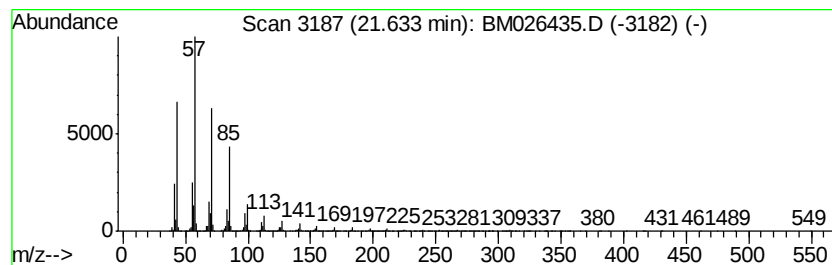
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Quant Title : SVOA CALIBRATION

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

Peak Number 27 (DEL) Alkane: Straight-Chai... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.63	21.55 ng/ul	2910640	Chrysene-d12	21.02

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heneicosane	296	C21H44	000629-94-7	95
2		Heptadecane, 9-hexyl-	324	C23H48	055124-79-3	94
3		Eicosane, 10-methyl-	296	C21H44	054833-23-7	93
4		Hexacosane	366	C26H54	000630-01-3	91
5		Heneicosane	296	C21H44	000629-94-7	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM061720\
Data File : BM026435.D
Acq On : 17 Jun 2020 20:43
Operator : JU/CG
Sample : L2959-08
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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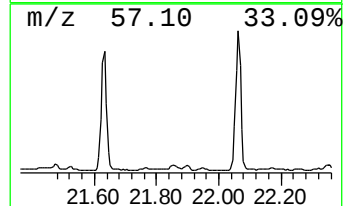
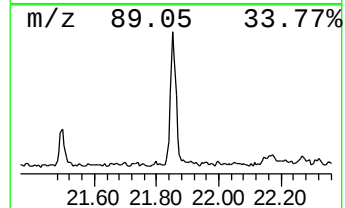
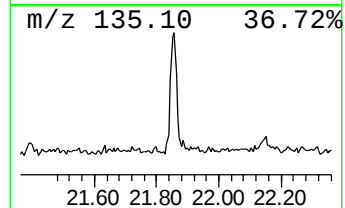
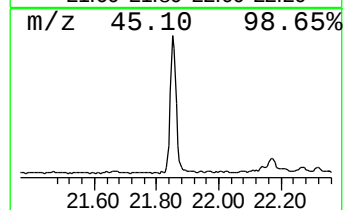
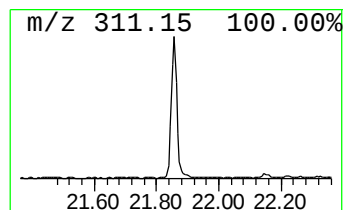
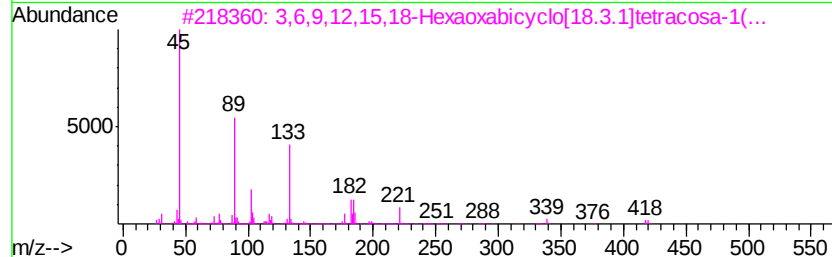
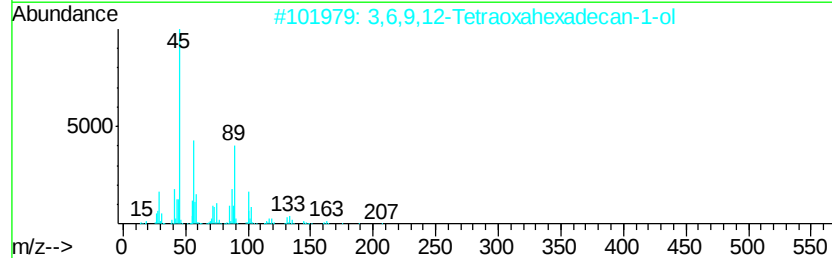
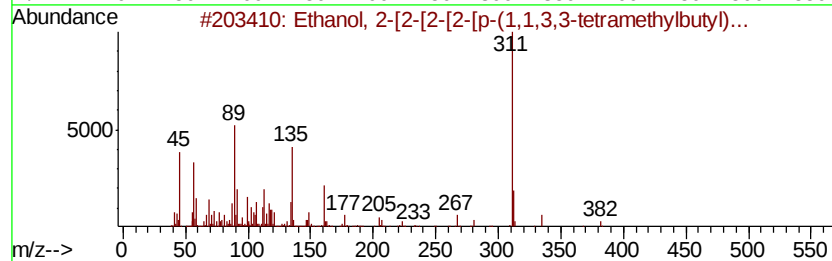
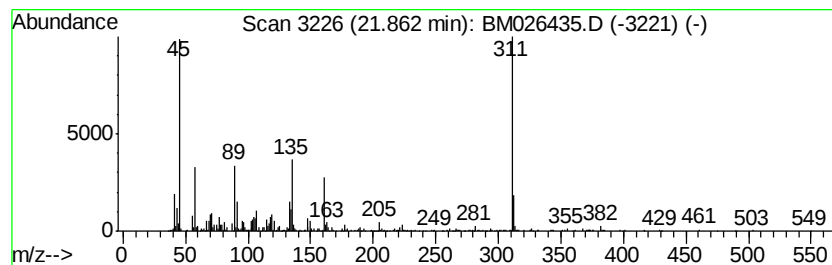
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Quant Title : SVOA CALIBRATION

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

Peak Number 28 unknown-02 Concentration Rank 29

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.86	3.41 ng/ul	459956	Chrysene-d12	21.02

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Ethanol, 2-[2-[2-[2-[p-(1,1,3,3-...	382	C22H38O5	002315-63-1	43
2			3,6,9,12-Tetraoxahexadecan-1-ol	250	C12H26O5	001559-34-8	27
3			3,6,9,12,15,18-Hexaoxabicyclo[18...	418	C18H27BrO6	111982-22-0	27
4			Hexaethylene glycol	282	C12H26O7	002615-15-8	27
5			3,6,9,12,15-Pentaoxabicyclo[15.3...	326	C17H26O6	071128-99-9	25



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM061720\
Data File : BM026435.D
Acq On : 17 Jun 2020 20:43
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Sample : L2959-08
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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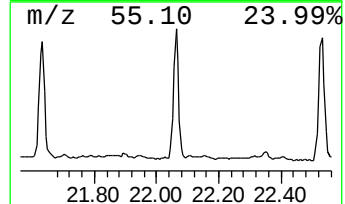
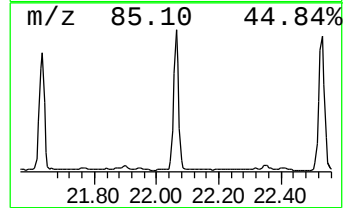
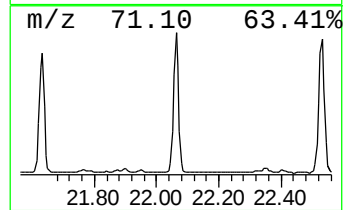
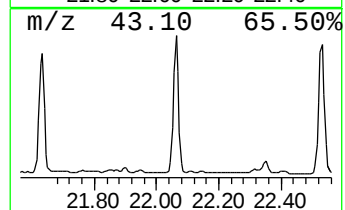
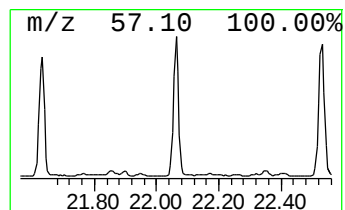
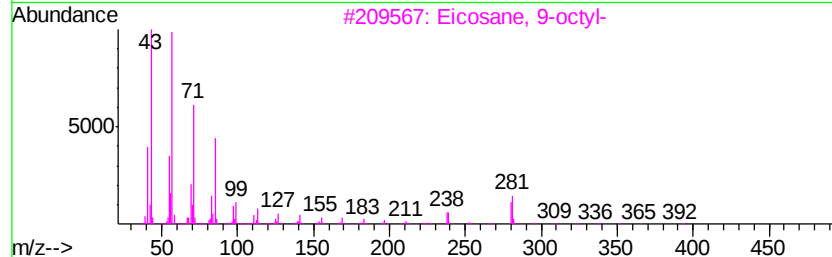
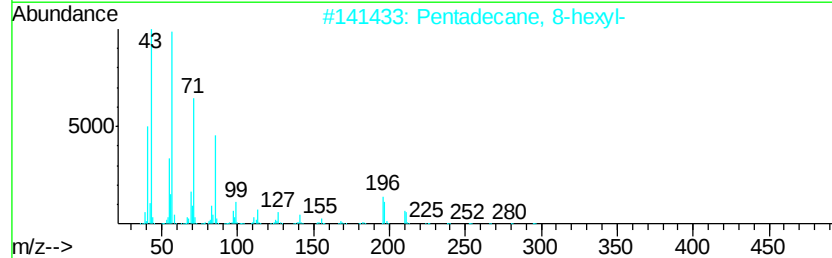
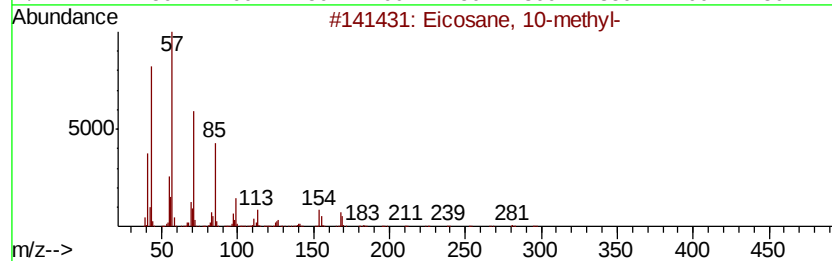
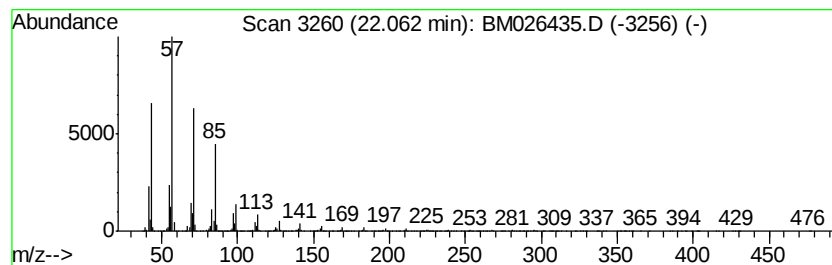
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Quant Title : SVOA CALIBRATION

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

Peak Number 29 (DEL) Alkane: Straight-Chai... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.06	24.71 ng/ul	3337290	Chrysene-d12	21.02

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Eicosane, 10-methyl-	296	C21H44	054833-23-7	94
2		Pentadecane, 8-hexyl-	296	C21H44	013475-75-7	94
3		Eicosane, 9-octyl-	394	C28H58	013475-77-9	94
4		Octadecane	254	C18H38	000593-45-3	93
5		2-methyloctacosane	408	C29H60	1000376-72-8	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM061720\
Data File : BM026435.D
Acq On : 17 Jun 2020 20:43
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Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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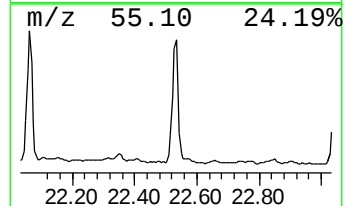
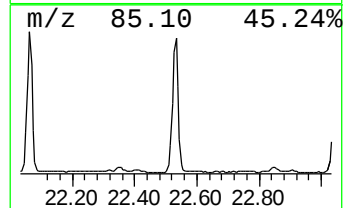
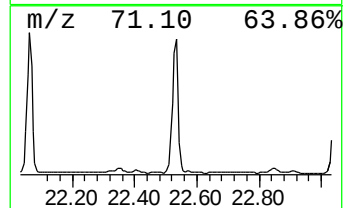
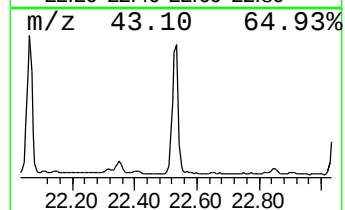
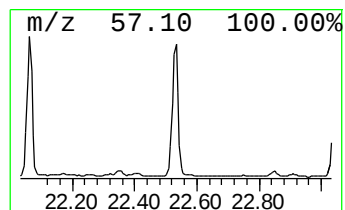
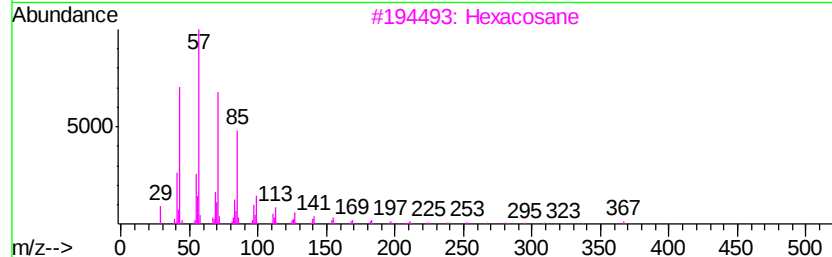
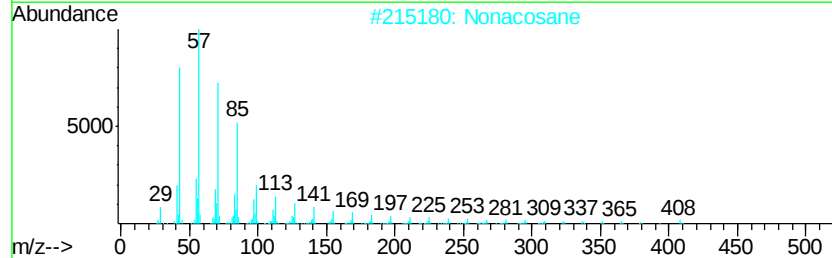
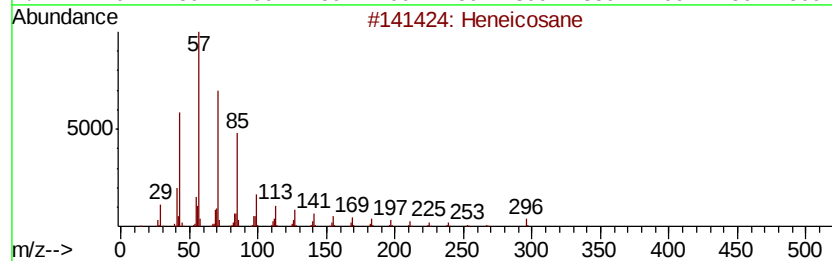
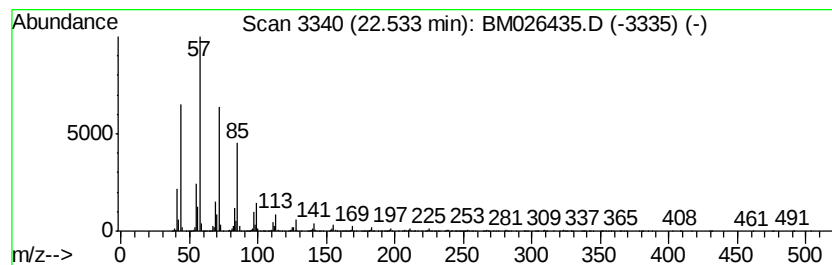
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Quant Title : SVOA CALIBRATION

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

Peak Number 30 (DEL) Alkane: Straight-Chai... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.53	23.71 ng/ul	3733500	Perylene-d12	23.11

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heneicosane	296	C21H44	000629-94-7	97
2		Nonacosane	408	C29H60	000630-03-5	95
3		Hexacosane	366	C26H54	000630-01-3	94
4		Eicosane, 10-methyl-	296	C21H44	054833-23-7	93
5		Heptacosane	380	C27H56	000593-49-7	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM061720\
Data File : BM026435.D
Acq On : 17 Jun 2020 20:43
Operator : JU/CG
Sample : L2959-08
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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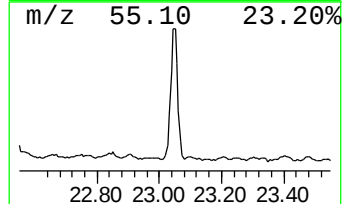
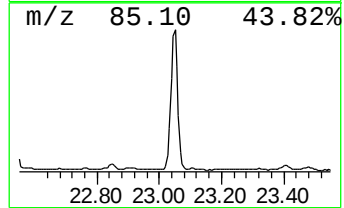
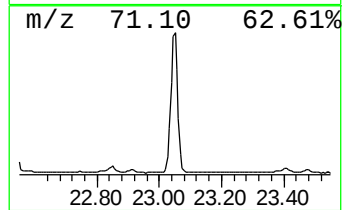
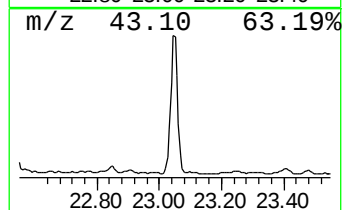
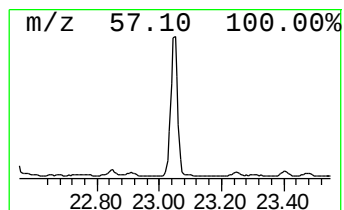
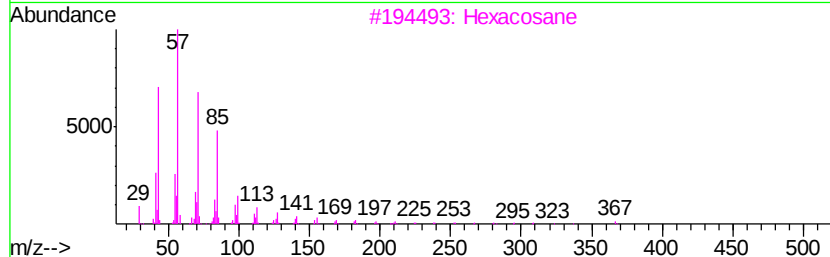
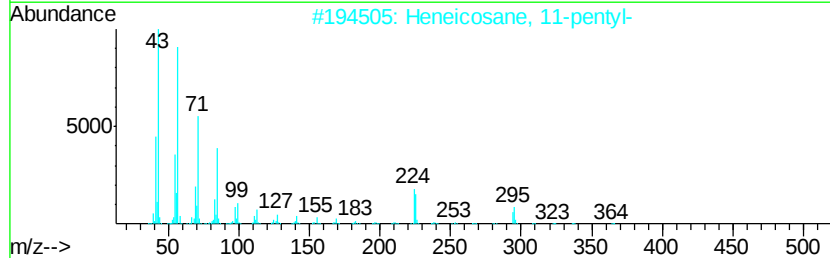
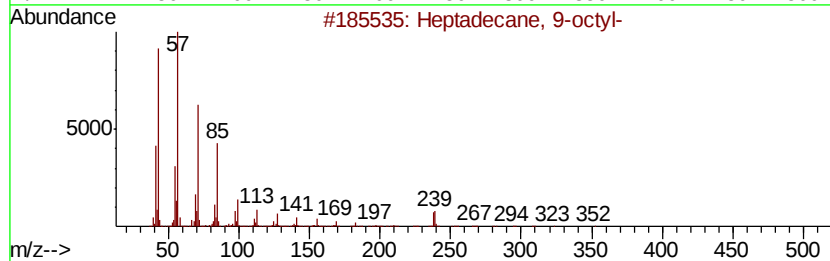
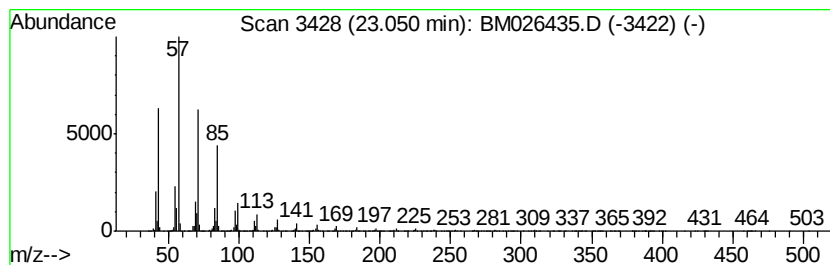
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Quant Title : SVOA CALIBRATION

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

Peak Number 31 (DEL) Alkane: Straight-Chai... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.05	23.91 ng/ul	3764410	Perylene-d12	23.11

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heptadecane, 9-octyl-	352	C25H52	007225-64-1	95
2		Heneicosane, 11-pentyl-	366	C26H54	014739-72-1	95
3		Hexacosane	366	C26H54	000630-01-3	94
4		Octadecane, 5,14-dibutyl-	366	C26H54	055282-13-8	94
5		Tetracosane	338	C24H50	000646-31-1	91



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Data File : BM026435.D
Acq On : 17 Jun 2020 20:43
Operator : JU/CG
Sample : L2959-08
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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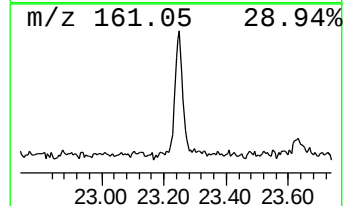
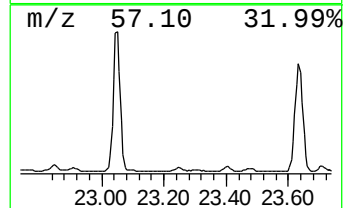
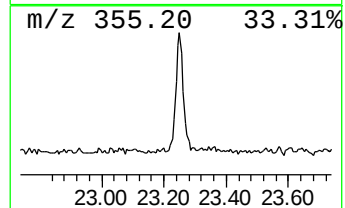
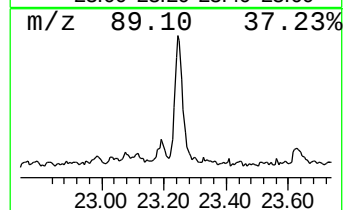
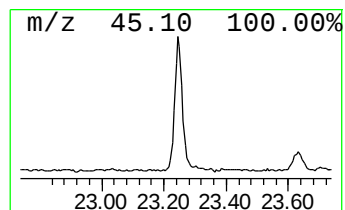
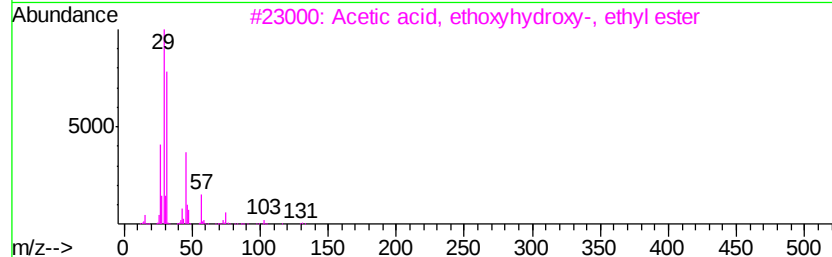
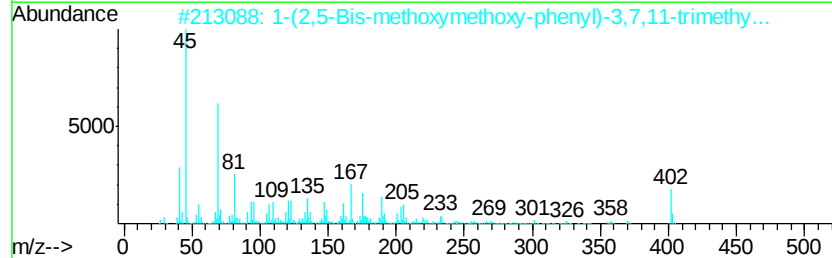
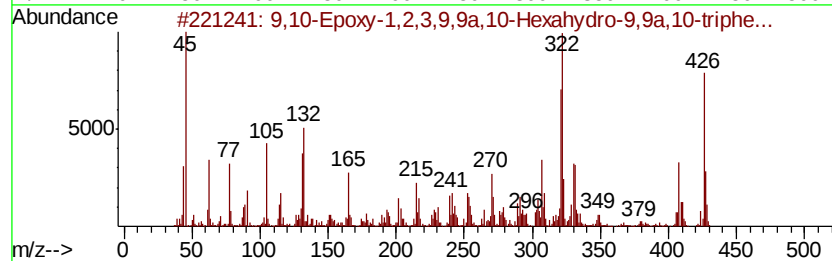
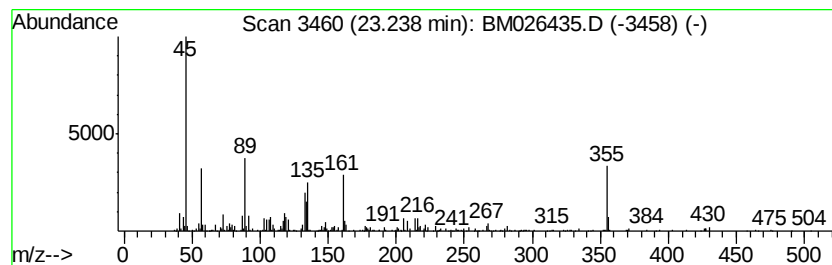
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Quant Title : SVOA CALIBRATION

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

Peak Number 33 unknown-03 Concentration Rank 30

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.24	3.40 ng/ul	535864	Perylene-d12	23.11

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	9,10-Epoxy-1,2,3,9,9a,10-Hexahyd...	426	C32H26O	1000326-57-1	30		
2	1-(2,5-Bis-methoxymethoxy-phenyl)...	402	C25H38O4	1000188-50-6	25		
3	Acetic acid, ethoxyhydroxy-, eth...	148	C6H12O4	049653-17-0	9		
4	Propanoic acid, 2-hydroxy-, meth...	104	C4H8O3	002155-30-8	9		
5	Diethyl carbonate	118	C5H10O3	000105-58-8	9		



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM061720\
Data File : BM026435.D
Acq On : 17 Jun 2020 20:43
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Sample : L2959-08
Misc :
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ClientSampleId :
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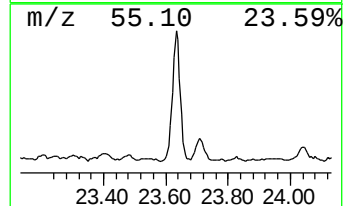
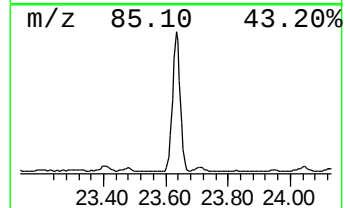
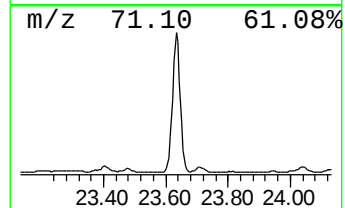
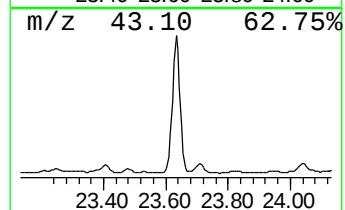
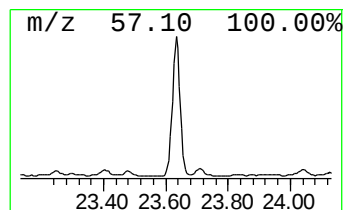
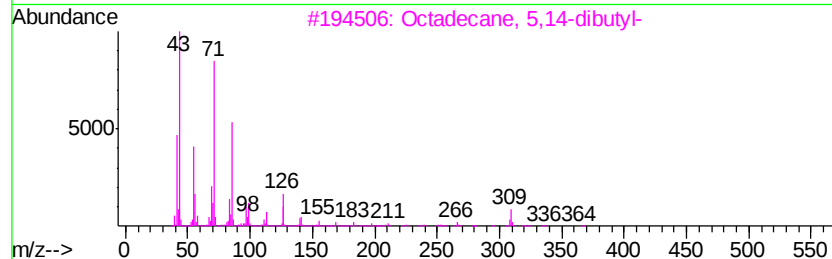
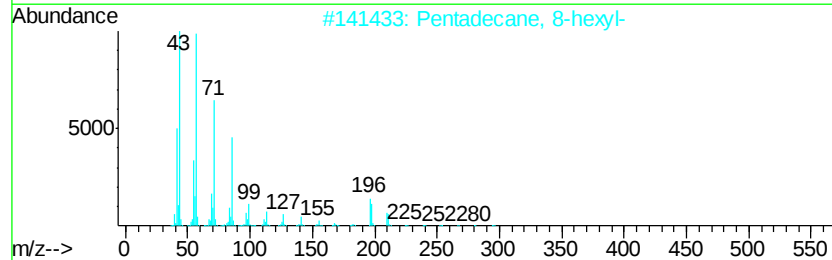
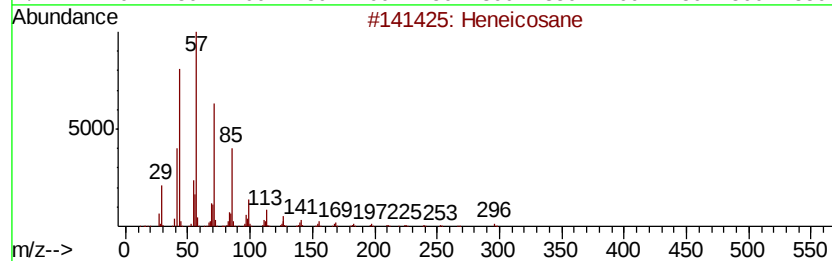
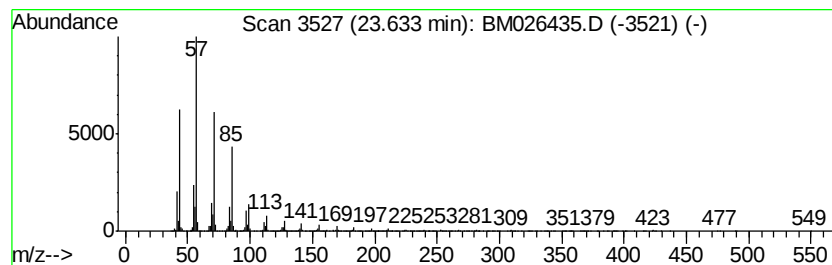
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Quant Title : SVOA CALIBRATION

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

Peak Number 34 (DEL) Alkane: Straight-Chai... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.63	18.95 ng/ul	2984030	Perylene-d12	23.11

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heneicosane	296	C21H44	000629-94-7	97
2		Pentadecane, 8-hexyl-	296	C21H44	013475-75-7	96
3		Octadecane, 5,14-dibutyl-	366	C26H54	055282-13-8	94
4		Octacosane	394	C28H58	000630-02-4	93
5		Tricosane	324	C23H48	000638-67-5	93



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM061720\
Data File : BM026435.D
Acq On : 17 Jun 2020 20:43
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Sample : L2959-08
Misc :
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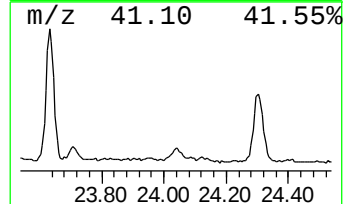
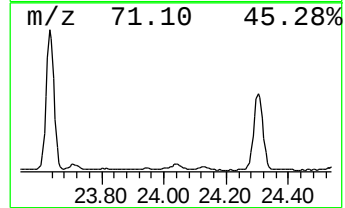
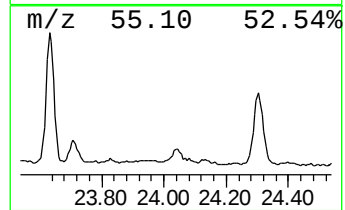
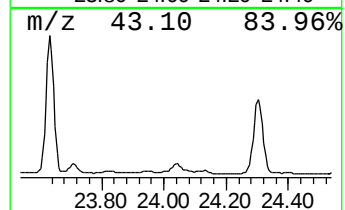
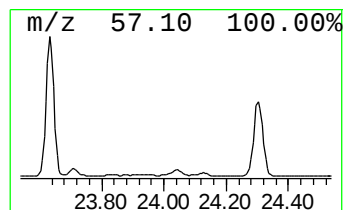
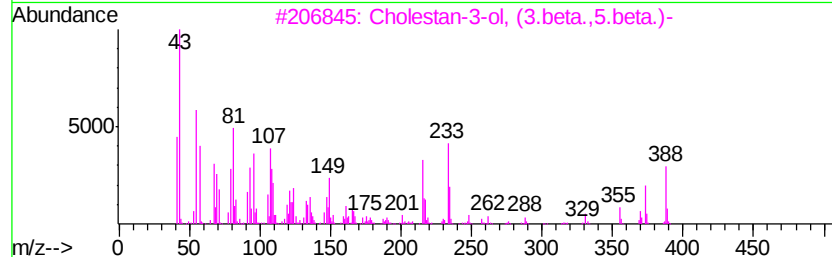
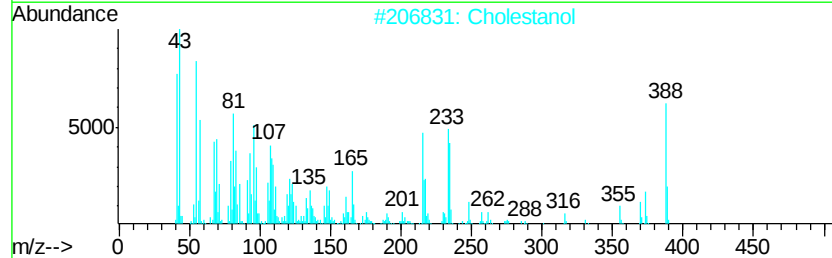
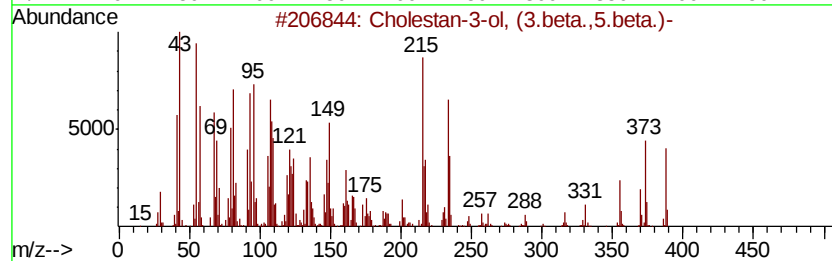
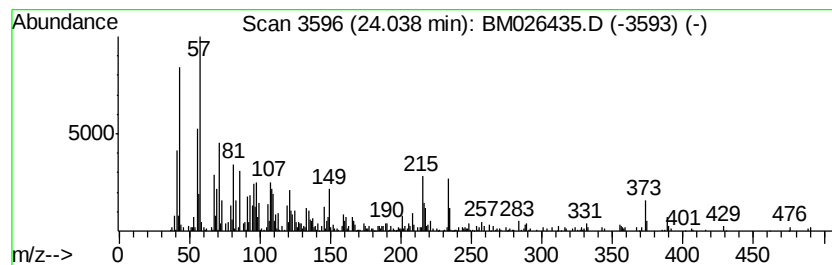
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Quant Title : SVOA CALIBRATION

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

Peak Number 35 Cholestan-3-ol, (3.beta.,5.... Concentration Rank 33

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.04	2.87 ng/ul	452222	Perylene-d12	23.11

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cholestan-3-ol, (3.beta.,5.beta.)-	388	C27H48O	000360-68-9	53
2		Cholestanol	388	C27H48O	000080-97-7	49
3		Cholestan-3-ol, (3.beta.,5.beta.)-	388	C27H48O	000360-68-9	38
4		Cholestan-3-ol	388	C27H48O	027409-41-2	35
5		Cholestan-3-ol	388	C27H48O	027409-41-2	12



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM061720\
Data File : BM026435.D
Acq On : 17 Jun 2020 20:43
Operator : JU/CG
Sample : L2959-08
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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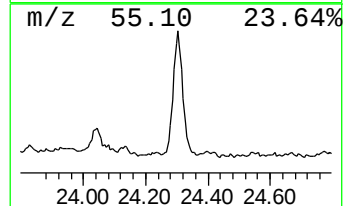
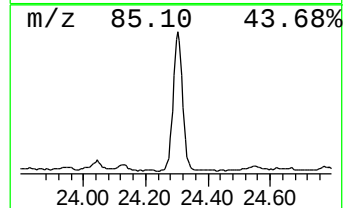
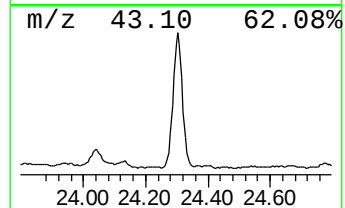
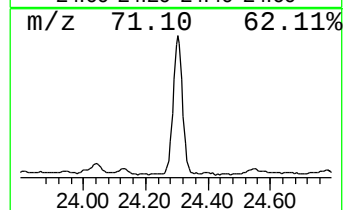
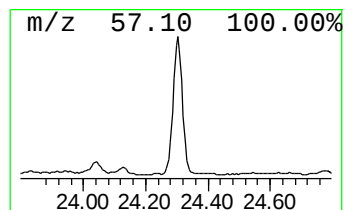
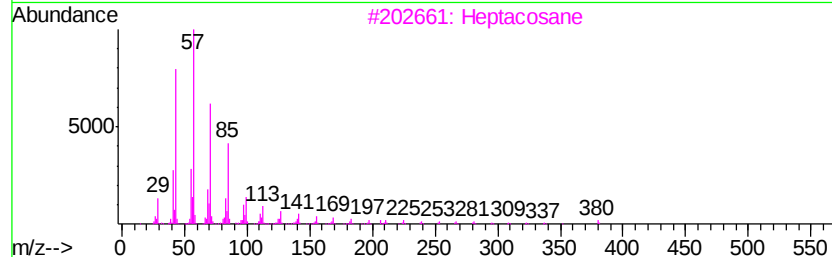
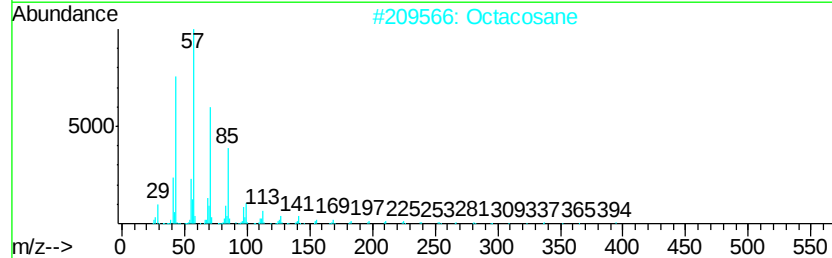
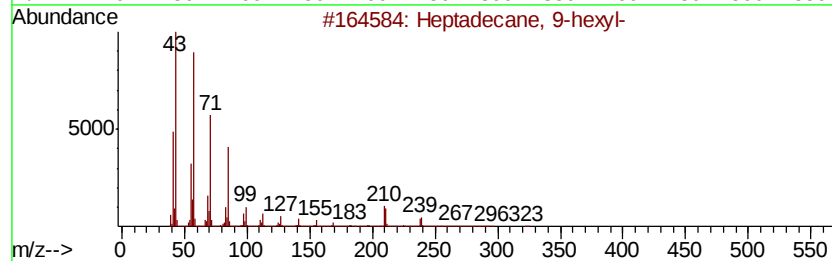
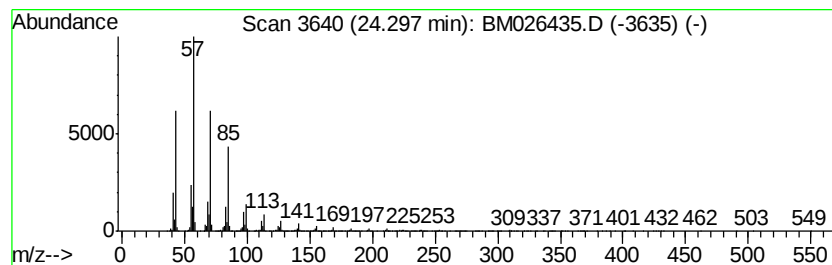
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Quant Title : SVOA CALIBRATION

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

Peak Number 36 (DEL) Alkane: Straight-Chai... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.30	12.86 ng/ul	2025280	Perylene-d12	23.11

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heptadecane, 9-hexyl-	324	C23H48	055124-79-3	94
2		Octacosane	394	C28H58	000630-02-4	91
3		Heptacosane	380	C27H56	000593-49-7	91
4		2-methyloctacosane	408	C29H60	1000376-72-8	91
5		Tetracosane, 11-decyl-	479	C34H70	055429-84-0	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM061720\
Data File : BM026435.D
Acq On : 17 Jun 2020 20:43
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Sample : L2959-08
Misc :
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ClientSampleId :
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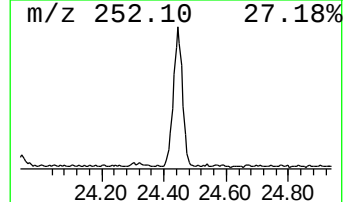
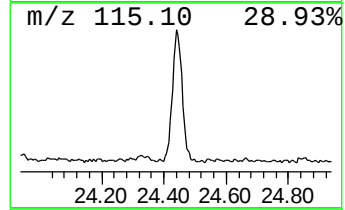
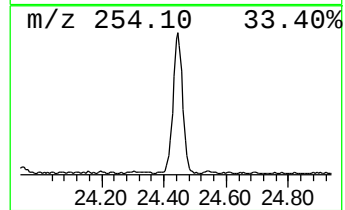
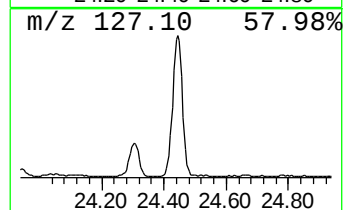
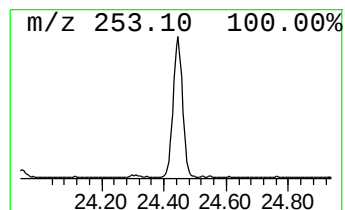
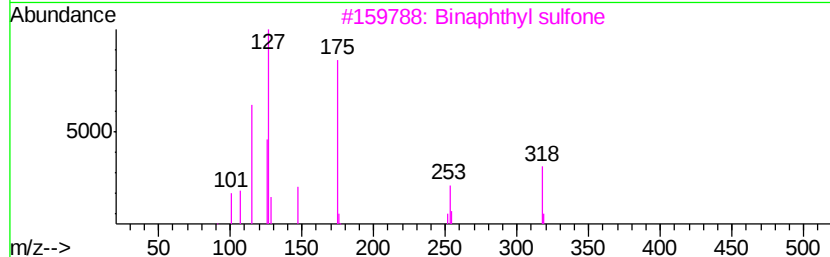
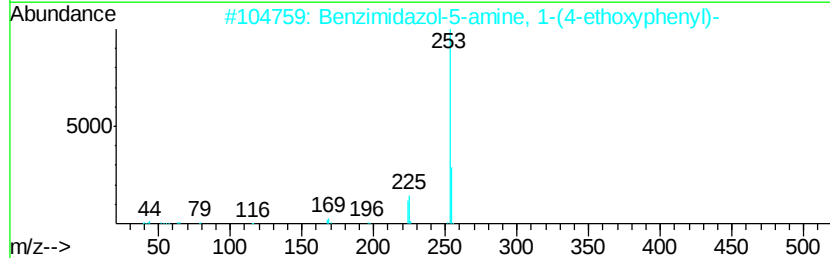
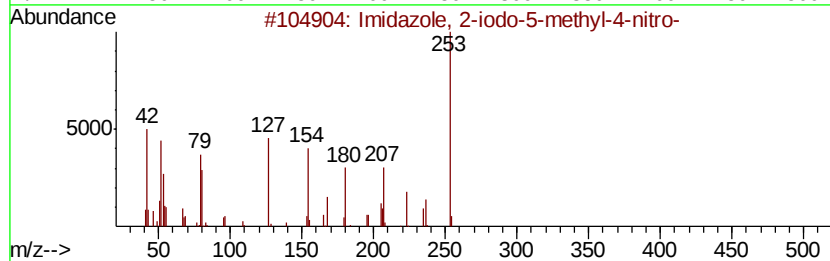
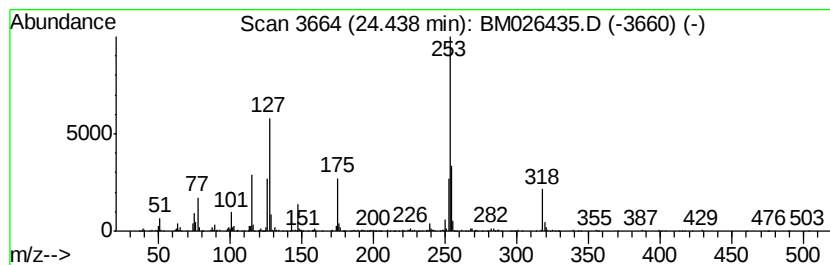
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Quant Title : SVOA CALIBRATION

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

Peak Number 37 unknown-04 Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.44	5.50 ng/ul	865723	Perylene-d12	23.11

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Imidazole, 2-iodo-5-methyl-4-nitro-	253	C4H4IN3O2	149510-86-1	38
2			Benzimidazol-5-amine, 1-(4-ethox...	253	C15H15N3O	007104-62-3	27
3			Binaphthyl sulfone	318	C20H14O2S	032390-26-4	22
4			2-Anilino-4-chloroquinoline	254	C15H11ClN2	032888-95-2	14
5			2-[2-Thienyl]-4-acetyl quinoline	253	C15H11NOS	027302-83-6	12



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM061720\
Data File : BM026435.D
Acq On : 17 Jun 2020 20:43
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Sample : L2959-08
Misc :
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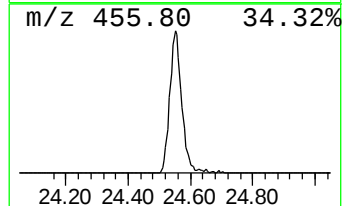
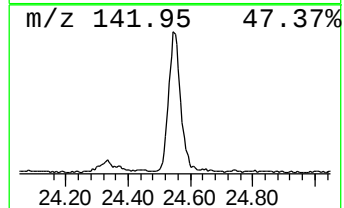
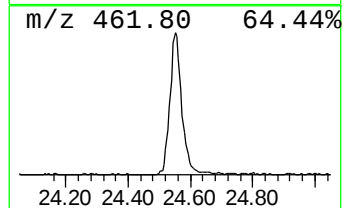
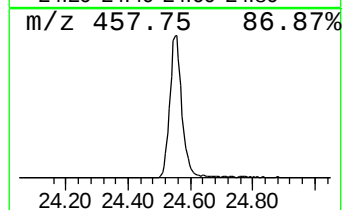
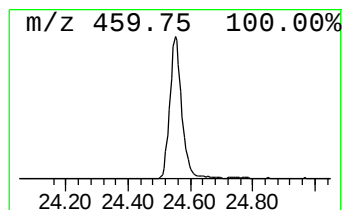
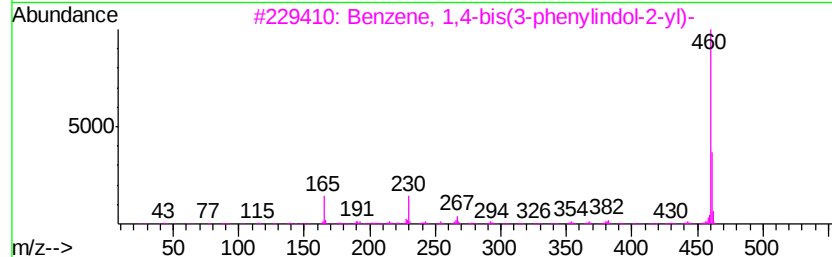
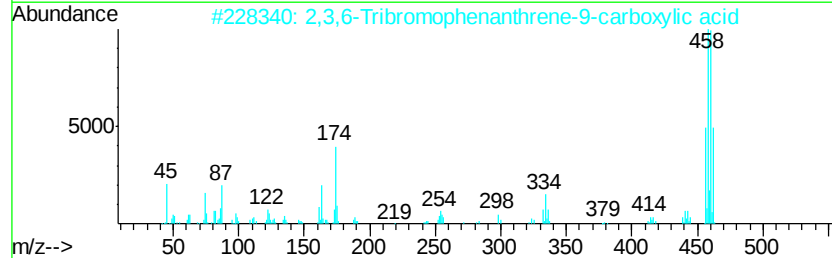
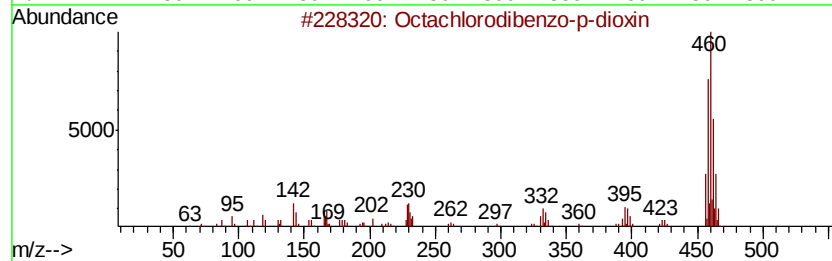
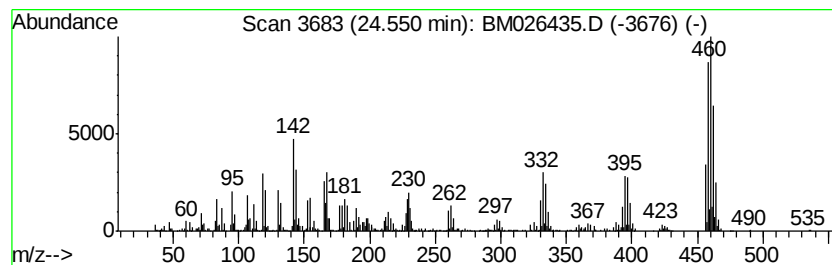
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Quant Title : SVOA CALIBRATION

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

Peak Number 38 Octachlorodibenzo-p-dioxin Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.55	11.26 ng/ul	1773300	Perylene-d12	23.11

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octachlorodibenzo-p-dioxin	456	C12Cl8O2	003268-87-9	83
2		2,3,6-Tribromophenanthrene-9-car...	456	C15H7Br3O2	056988-60-4	12
3		Benzene, 1,4-bis(3-phenylindol-2...	460	C34H24N2	164936-88-3	10
4		(1-Amino-5-thiophen-2-yl-6,7,8,9...	458	C22H16Cl2N2OS2	1000318-00-6	9
5		Pyridine, pentaphenyl-	459	C35H25N	040249-26-1	9



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM061720\
Data File : BM026435.D
Acq On : 17 Jun 2020 20:43
Operator : JU/CG
Sample : L2959-08
Misc :
ALS Vial : 11 Sample Multiplier: 1

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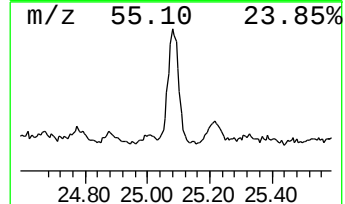
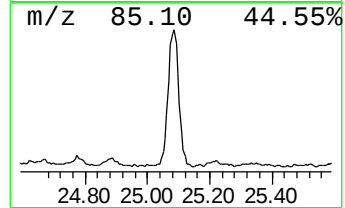
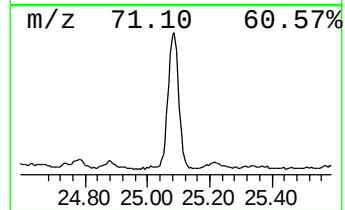
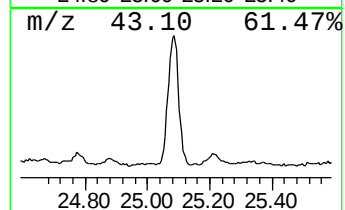
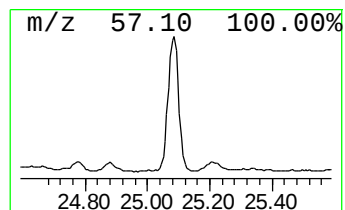
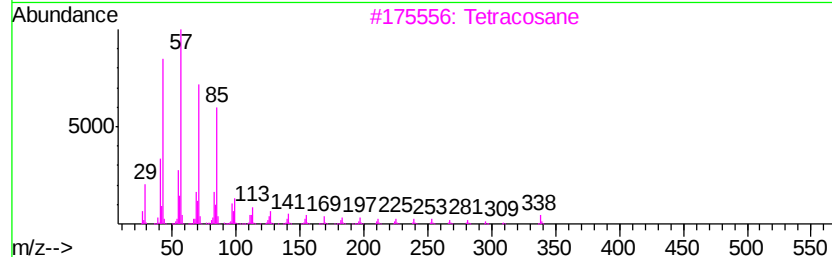
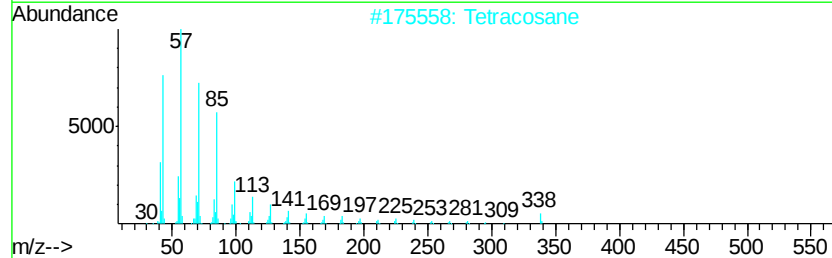
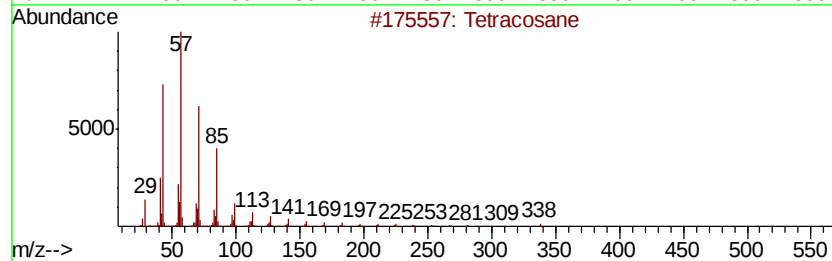
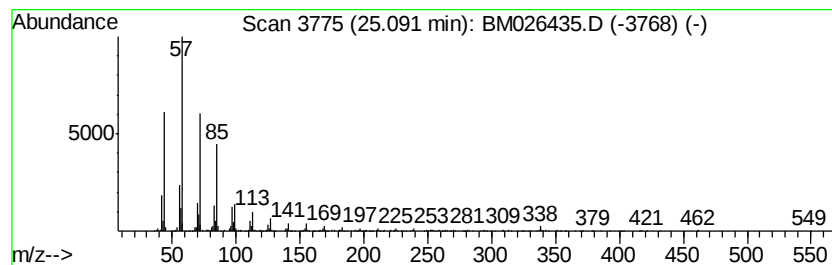
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Quant Title : SVOA CALIBRATION

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

Peak Number 39 (DEL) Alkane: Straight-Chai... Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
25.09	5.66 ng/ul	891005	Perylene-d12	23.11

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tetracosane	338	C24H50	000646-31-1	98
2			Tetracosane	338	C24H50	000646-31-1	97
3			Tetracosane	338	C24H50	000646-31-1	96
4			Tetracosane	338	C24H50	000646-31-1	96
5			Hexacosane	366	C26H54	000630-01-3	94



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM061720\
Data File : BM026435.D
Acq On : 17 Jun 2020 20:43
Operator : JU/CG
Sample : L2959-08
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
BG2K2

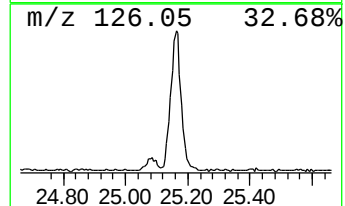
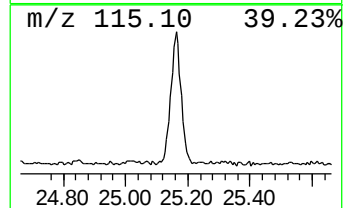
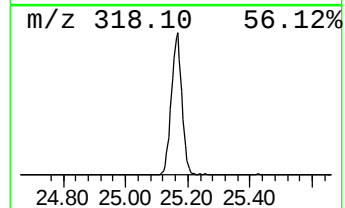
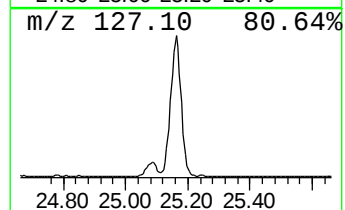
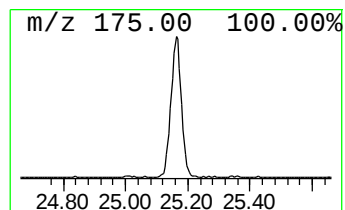
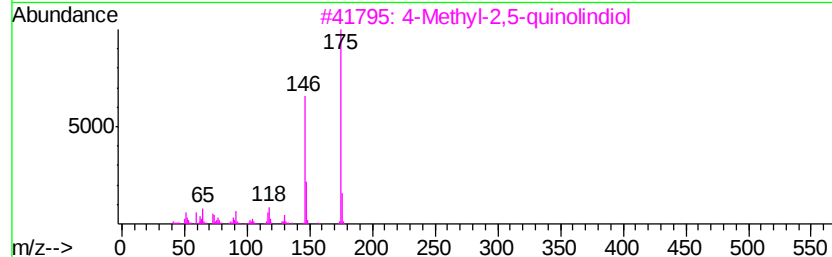
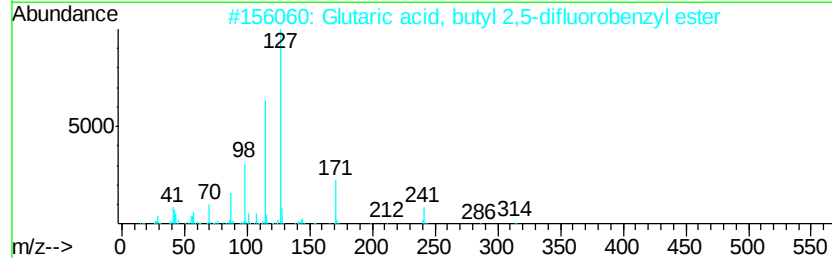
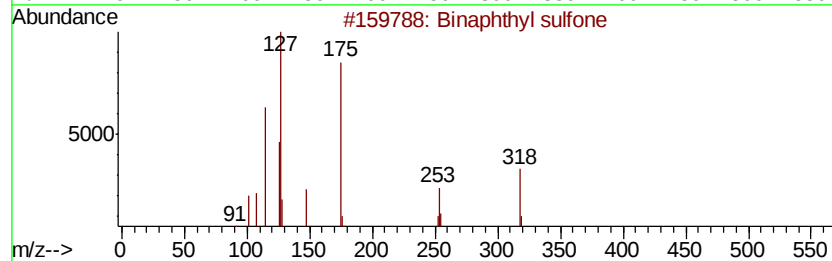
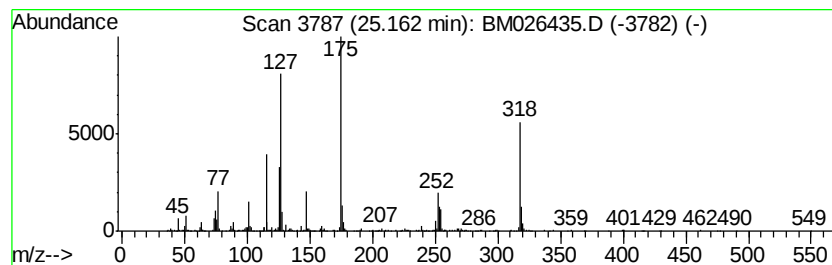
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Quant Title : SVOA CALIBRATION

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

Peak Number 40 unknown-05 Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
25.16	5.65 ng/ul	890264	Perylene-d12	23.11

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Binaphthyl sulfone	318	C20H14O2S	032390-26-4	43
2			Glutaric acid, butyl 2,5-difluor...	314	C16H20F2O4	1000376-94-5	22
3			4-Methyl-2,5-quinolindiol	175	C10H9NO2	131195-67-0	14
4			2,2'-Dithiodi[3-pyridinol]	252	C10H8N2O2S2	068750-09-4	14
5			4-Methyl-2,6-dihydroxyquinoline	175	C10H9NO2	034982-01-9	14



Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM061720\
 Data File : BM026435.D
 Acq On : 17 Jun 2020 20:43
 Operator : JU/CG
 Sample : L2959-08
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 BG2K2

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM061220MA.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...	7.04	4.4	ng/ul	210889	1	7.39	957446	20.0
D-Limonene	7.62	7.1	ng/ul	339498	1	7.39	957446	20.0
(DEL) Alkane: Str...	7.93	2.2	ng/ul	104944	1	7.39	957446	20.0
(DEL) Alkane: Str...	8.63	4.9	ng/ul	235276	1	7.39	957446	20.0
(DEL) Alkane: Cyc...	8.73	2.5	ng/ul	119871	1	7.39	957446	20.0
1,2,3-Propanetriol...	10.70	2.4	ng/ul	166571	2	10.15	1393490	20.0
Phenol, 3,5-dichl...	12.77	2.0	ng/ul	197340	3	14.05	1932700	20.0
Phenol, 3,4-dichl...	13.09	2.9	ng/ul	279462	3	14.05	1932700	20.0
Dodecanoic acid	14.53	5.1	ng/ul	495340	3	14.05	1932700	20.0
Tributyl phosphate	15.46	8.8	ng/ul	1008530	4	16.82	2297130	20.0
Ethanol, 2-[4-(1,...	17.19	3.9	ng/ul	451018	4	16.82	2297130	20.0
(DEL) Alkane: Str...	17.34	3.0	ng/ul	350710	4	16.82	2297130	20.0
[1,1'-Biphenyl]-2...	17.43	2.2	ng/ul	254264	4	16.82	2297130	20.0
1-Iodoundecane	18.04	3.5	ng/ul	396728	4	16.82	2297130	20.0
(DEL) Alkane: Str...	18.68	6.6	ng/ul	754187	4	16.82	2297130	20.0
trans-13-Octadec...	18.93	22.1	ng/ul	2983400	5	21.02	2701590	20.0
Octadecanoic acid	19.06	14.5	ng/ul	1956050	5	21.02	2701590	20.0
Ethanol, 2-[2-[4-...	19.12	4.7	ng/ul	637047	5	21.02	2701590	20.0
(DEL) Alkane: Str...	19.27	8.6	ng/ul	1154830	5	21.02	2701590	20.0
(DEL) Alkane: Str...	19.81	10.4	ng/ul	1404850	5	21.02	2701590	20.0
(DEL) Alkane: Str...	20.31	12.7	ng/ul	1721090	5	21.02	2701590	20.0
Ethanol, 2-[2-[2-...	20.62	4.1	ng/ul	559610	5	21.02	2701590	20.0
unknown-01	20.73	2.1	ng/ul	286364	5	21.02	2701590	20.0
1-Octadecanesulph...	20.77	20.8	ng/ul	2806570	5	21.02	2701590	20.0
(DEL) Alkane: Str...	21.22	16.9	ng/ul	2289650	5	21.02	2701590	20.0
(DEL) Alkane: Str...	21.63	21.6	ng/ul	2910640	5	21.02	2701590	20.0
unknown-02	21.86	3.4	ng/ul	459956	5	21.02	2701590	20.0
(DEL) Alkane: Str...	22.06	24.7	ng/ul	3337290	5	21.02	2701590	20.0
(DEL) Alkane: Str...	22.53	23.7	ng/ul	3733500	6	23.11	3149010	20.0
(DEL) Alkane: Str...	23.05	23.9	ng/ul	3764410	6	23.11	3149010	20.0
1,2,3,7,8,9-Hexac...	23.20	2.6	ng/ul	405341	6	23.11	3149010	20.0
unknown-03	23.24	3.4	ng/ul	535864	6	23.11	3149010	20.0
(DEL) Alkane: Str...	23.63	18.9	ng/ul	2984030	6	23.11	3149010	20.0
Cholestan-3-ol, (...	24.04	2.9	ng/ul	452222	6	23.11	3149010	20.0
(DEL) Alkane: Str...	24.30	12.9	ng/ul	2025280	6	23.11	3149010	20.0
unknown-04	24.44	5.5	ng/ul	865723	6	23.11	3149010	20.0
Octachlorodibenzo...	24.55	11.3	ng/ul	1773300	6	23.11	3149010	20.0
(DEL) Alkane: Str...	25.09	5.7	ng/ul	891005	6	23.11	3149010	20.0
unknown-05	25.16	5.7	ng/ul	890264	6	23.11	3149010	20.0