

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN022220\
 Data File : BN009850.D
 Acq On : 22 Feb 2020 17:45
 Operator : CG/JU
 Sample : L1516-12
 Misc :
 ALS Vial : 39 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 DBD30

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_N\METHODS\SOM-EPA-BN021220MA.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	3.034	4	8	16	rBV	40587	55953	1.92%	0.146%
2	3.375	62	66	68	rBV4	3815	6517	0.22%	0.017%
3	3.722	121	125	131	rVB3	5737	8925	0.31%	0.023%
4	4.328	224	228	233	rVB6	2621	5254	0.18%	0.014%
5	4.599	270	274	282	rBV7	9704	22207	0.76%	0.058%
6	6.593	606	613	624	rBV	659235	1100153	37.74%	2.862%
7	6.746	633	639	651	rBV	461791	742256	25.46%	1.931%
8	6.934	664	671	684	rBV	629094	1046261	35.89%	2.722%
9	7.240	721	723	725	rBV3	3118	3503	0.12%	0.009%
10	7.387	742	748	758	rBV	504103	794364	27.25%	2.067%
11	7.475	760	763	766	rVV3	11883	14653	0.50%	0.038%
12	8.104	863	870	883	rBV	730484	1230686	42.22%	3.202%
13	8.216	886	889	895	rVB4	3090	4267	0.15%	0.011%
14	8.534	936	943	955	rBV	637529	1070030	36.71%	2.784%
15	8.722	971	975	977	rVB5	3289	4677	0.16%	0.012%
16	8.975	1012	1018	1022	rBV	14481	18271	0.63%	0.048%
17	9.246	1057	1064	1082	rBV	528132	895807	30.73%	2.331%
18	9.604	1122	1125	1128	rBV2	2584	3436	0.12%	0.009%
19	9.775	1148	1154	1172	rBV	693413	1207819	41.44%	3.142%
20	10.134	1208	1215	1225	rBV	674492	1109229	38.05%	2.886%
21	10.287	1235	1241	1262	rBV	594080	1226604	42.08%	3.191%
22	11.010	1359	1364	1369	rBV4	2161	4596	0.16%	0.012%
23	11.393	1426	1429	1432	rBV3	3358	4388	0.15%	0.011%
24	11.757	1486	1491	1494	rBV3	6233	8533	0.29%	0.022%
25	12.945	1685	1693	1700	rBV	284187	436269	14.97%	1.135%
26	13.269	1744	1748	1756	rBV4	5989	8433	0.29%	0.022%
27	13.457	1774	1780	1785	rBV	1173566	1677830	57.56%	4.365%
28	13.504	1785	1788	1796	rVV	124000	177987	6.11%	0.463%
29	13.575	1798	1800	1806	rVB5	3280	4226	0.14%	0.011%
30	13.716	1818	1824	1834	rBV	1352598	1985334	68.11%	5.165%
31	13.969	1862	1867	1871	rVB5	3515	5021	0.17%	0.013%
32	14.028	1871	1877	1890	rVV2	974930	1405119	48.21%	3.656%
33	14.275	1912	1919	1940	rBV	376995	847593	29.08%	2.205%
34	14.410	1940	1942	1946	rVB4	4668	6053	0.21%	0.016%

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Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN021220MA.M
 Title : SVOA CALIBRATION

35	14.486	1953	1955	1959	rVB3	8273	9316	0.32%	0.024%
36	14.869	2015	2020	2023	rBV4	3649	6104	0.21%	0.016%
37	14.928	2025	2030	2033	rBV3	7820	10059	0.35%	0.026%
38	15.034	2042	2048	2059	rBV	1750813	2442706	83.80%	6.355%
39	15.192	2069	2075	2086	rBV	719155	1111017	38.12%	2.890%
40	15.281	2086	2090	2095	rVV	22057	35922	1.23%	0.093%
41	15.428	2113	2115	2120	rVB6	2477	3983	0.14%	0.010%
42	15.492	2123	2126	2130	rVB5	2932	4008	0.14%	0.010%
43	15.792	2172	2177	2179	rBV4	3507	4303	0.15%	0.011%
44	15.822	2179	2182	2186	rVB4	5507	6456	0.22%	0.017%
45	16.039	2214	2219	2221	rBV3	3121	4170	0.14%	0.011%
46	16.586	2306	2312	2316	rBV6	5911	11550	0.40%	0.030%
47	16.786	2340	2346	2353	rBV	1218718	1647234	56.51%	4.285%
48	16.886	2357	2363	2371	rBV	1831315	2521849	86.52%	6.561%
49	17.128	2397	2404	2408	rBV2	23320	33061	1.13%	0.086%
50	17.204	2413	2417	2424	rBV	45632	61284	2.10%	0.159%
51	17.316	2432	2436	2438	rBV5	3050	4302	0.15%	0.011%
52	17.498	2465	2467	2472	rVB5	4367	5198	0.18%	0.014%
53	17.592	2479	2483	2488	rBV6	4527	10174	0.35%	0.026%
54	17.657	2490	2494	2496	rVV4	3770	4718	0.16%	0.012%
55	17.716	2498	2504	2518	rVV	264190	422330	14.49%	1.099%
56	17.828	2521	2523	2530	rVV8	9800	23648	0.81%	0.062%
57	17.880	2530	2532	2535	rVV3	8081	10797	0.37%	0.028%
58	17.910	2535	2537	2546	rVV8	6745	15366	0.53%	0.040%
59	18.004	2549	2553	2555	rVV2	9615	14283	0.49%	0.037%
60	18.039	2557	2559	2563	rVB5	3379	4320	0.15%	0.011%
61	18.145	2575	2577	2583	rBV6	5781	10408	0.36%	0.027%
62	18.192	2583	2585	2589	rVB5	3301	3507	0.12%	0.009%
63	18.398	2618	2620	2624	rVB4	4226	3392	0.12%	0.009%
64	18.533	2637	2643	2644	rBV5	5947	9494	0.33%	0.025%
65	18.586	2647	2652	2661	rBV	141177	193342	6.63%	0.503%
66	18.657	2662	2664	2666	rBV3	5322	3720	0.13%	0.010%
67	18.722	2671	2675	2678	rBV5	5510	7879	0.27%	0.020%
68	18.827	2690	2693	2696	rBV	14861	19052	0.65%	0.050%
69	18.863	2696	2699	2700	rVV3	10291	8568	0.29%	0.022%
70	18.910	2704	2707	2712	rVB	42147	53231	1.83%	0.138%
71	19.010	2720	2724	2732	rBV	74012	124584	4.27%	0.324%

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72	19.080	2733	2736	2740	rVB2	15869	23170	0.79%	0.060%
73	19.192	2748	2755	2762	rBV	2182890	2914818	100.00%	7.583%
74	19.286	2767	2771	2773	rVV5	6831	10692	0.37%	0.028%
75	19.357	2781	2783	2785	rVB4	3915	3363	0.12%	0.009%
76	19.427	2793	2795	2796	rBV2	5858	4783	0.16%	0.012%
77	19.522	2809	2811	2813	rBV3	5834	5082	0.17%	0.013%
78	19.751	2847	2850	2853	rBV3	19463	23435	0.80%	0.061%
79	19.786	2853	2856	2859	rVB3	15980	17209	0.59%	0.045%
80	19.857	2867	2868	2873	rBV5	5817	6161	0.21%	0.016%
81	20.116	2908	2912	2915	rBV4	22912	27875	0.96%	0.073%
82	20.163	2917	2920	2925	rVB6	11990	18932	0.65%	0.049%
83	20.286	2939	2941	2946	rVB3	23698	23268	0.80%	0.061%
84	20.704	3007	3012	3015	rBV	43327	73461	2.52%	0.191%
85	20.745	3015	3019	3026	rVV2	664184	863419	29.62%	2.246%
86	20.804	3026	3029	3038	rVB	164999	232833	7.99%	0.606%
87	20.951	3051	3054	3058	rBV	46427	50868	1.75%	0.132%
88	21.004	3058	3063	3075	rBV	1372792	1796668	61.64%	4.674%
89	21.133	3083	3085	3090	rBV5	18722	24874	0.85%	0.065%
90	21.198	3092	3096	3100	rVV	70134	82825	2.84%	0.215%
91	21.616	3163	3167	3173	rBV2	116043	157421	5.40%	0.410%
92	22.045	3236	3240	3244	rBV	161992	184742	6.34%	0.481%
93	22.157	3257	3259	3263	rVB	45114	45849	1.57%	0.119%
94	22.510	3316	3319	3324	rBV	167354	218465	7.49%	0.568%
95	22.974	3392	3398	3404	rBV	1656732	2805344	96.24%	7.298%
96	23.027	3404	3407	3412	rVV	207751	293302	10.06%	0.763%
97	23.104	3414	3420	3429	rVB	1008595	1782341	61.15%	4.637%
98	23.610	3502	3506	3513	rVB	164530	272839	9.36%	0.710%
99	23.686	3515	3519	3526	rVB2	125259	216302	7.42%	0.563%
100	24.274	3615	3619	3626	rVB	123092	250397	8.59%	0.651%

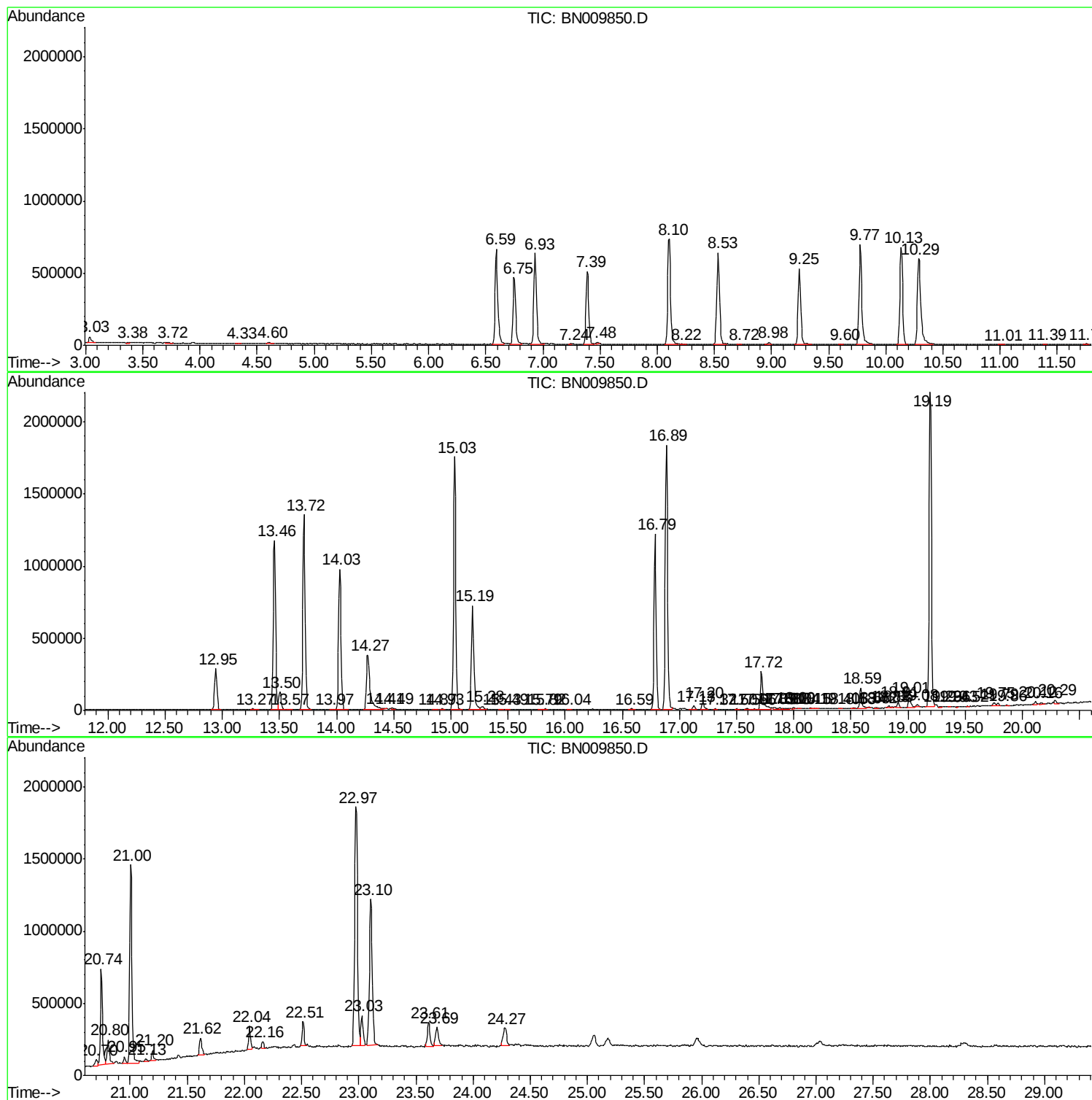
Sum of corrected areas: 38438327

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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN021220MA.M
Quant Title : SVOA CALIBRATION

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



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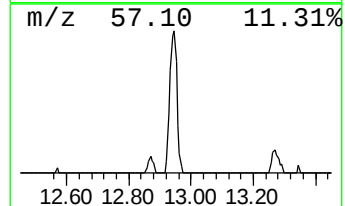
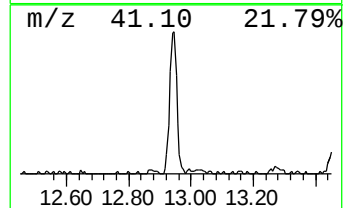
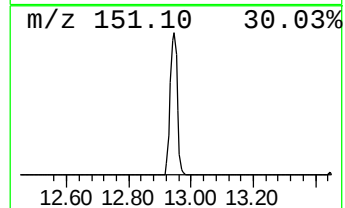
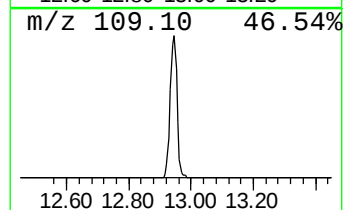
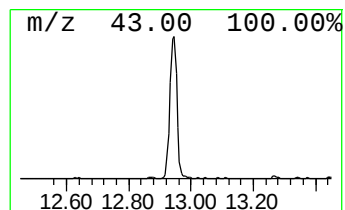
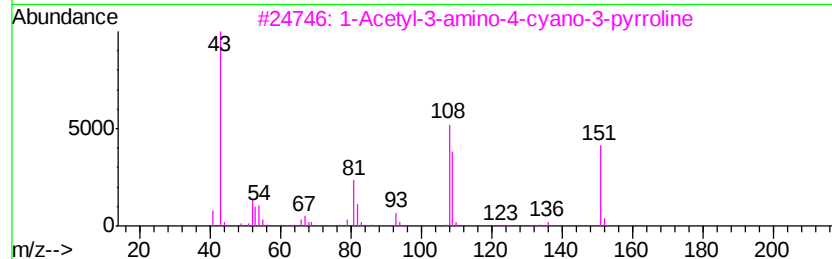
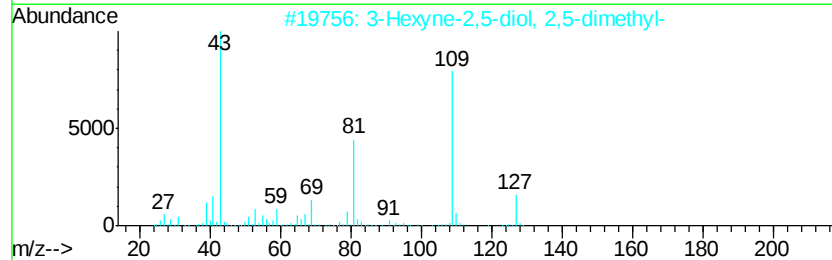
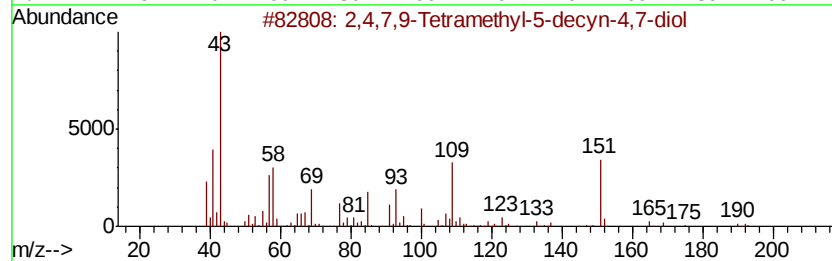
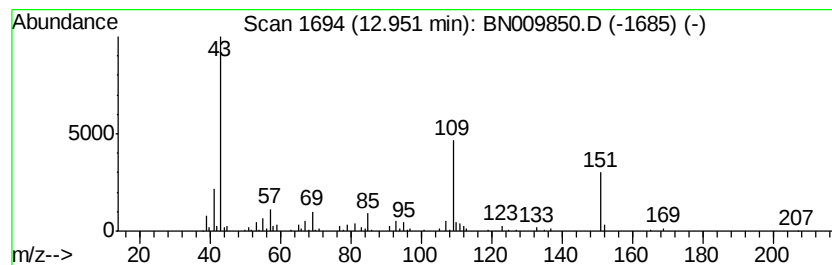
Instrument :
BNA_N
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN021220MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 unknown-01 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.	
12.95	6.21 ng/ul	436269	Acenaphthene-d10	14.03	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,4,7,9-Tetramethyl-5-decyn-4,7-...	226	C14H26O2	000126-86-3	32
2	3-Hexyne-2,5-diol, 2,5-dimethyl-	142	C8H14O2	000142-30-3	23
3	1-Acetyl-3-amino-4-cyano-3-pyrro...	151	C7H9N3O	002125-74-8	9
4	3-Hexyne-2,5-diol, 2,5-dimethyl-	142	C8H14O2	000142-30-3	9
5	Oxalic acid, allyl isohexyl ester	214	C11H18O4	1000309-23-3	9



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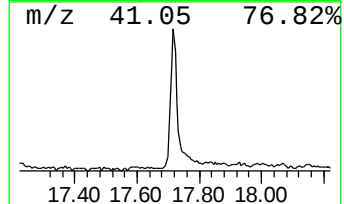
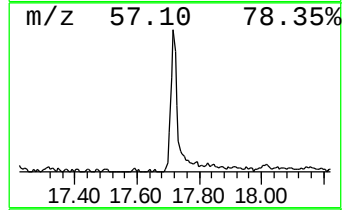
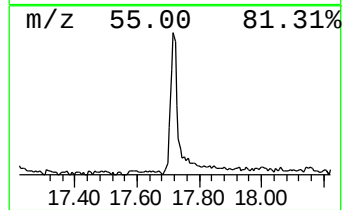
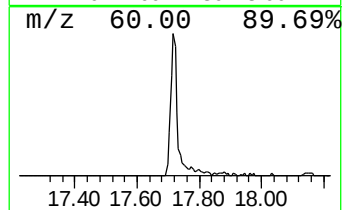
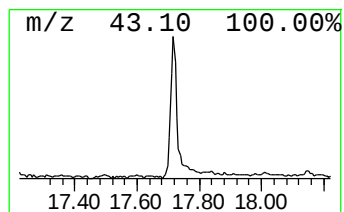
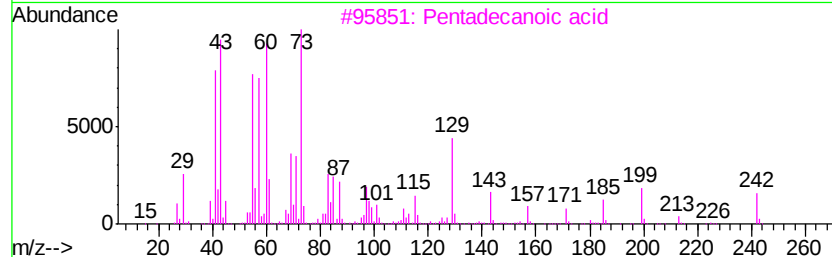
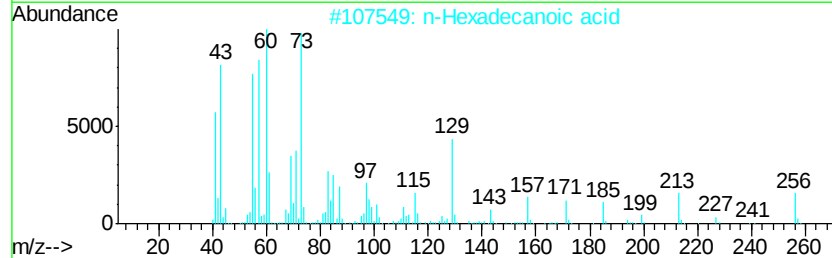
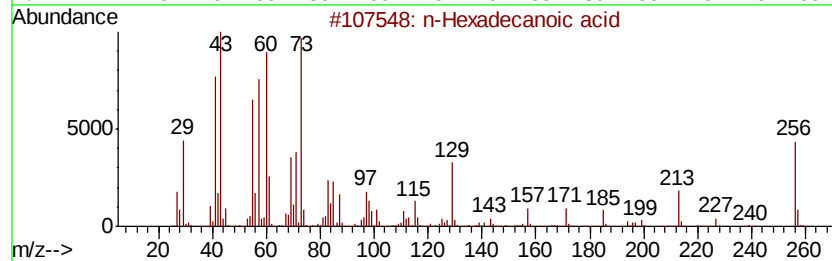
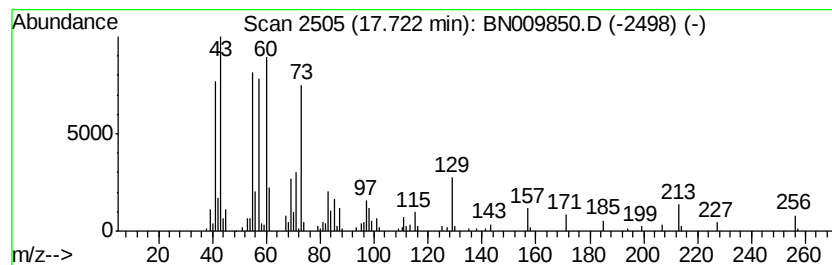
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN021220MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 n-Hexadecanoic acid Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.72	5.13 ng/ul	422330	Phenanthrene-d10	16.79

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	96
2			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	92
3			Pentadecanoic acid	242	C15H30O2	001002-84-2	90
4			Pentadecanoic acid	242	C15H30O2	001002-84-2	86
5			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	83



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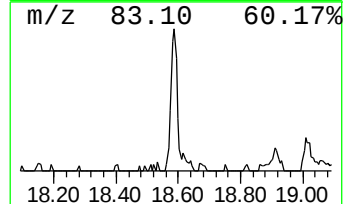
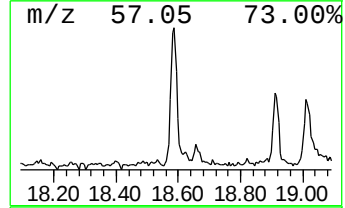
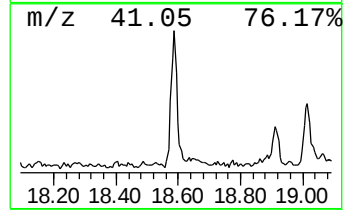
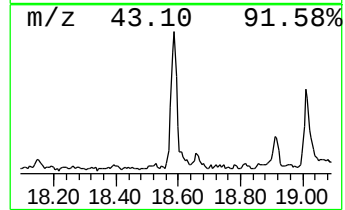
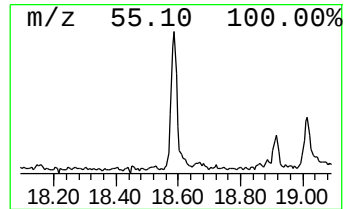
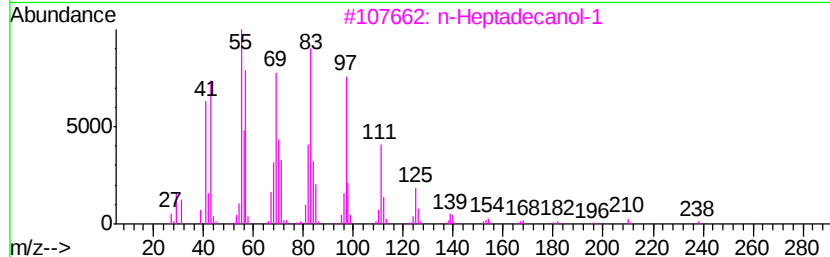
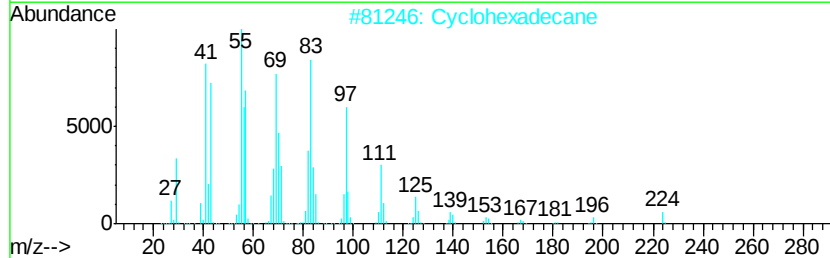
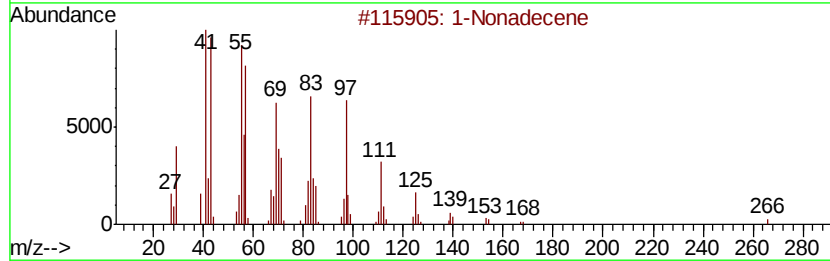
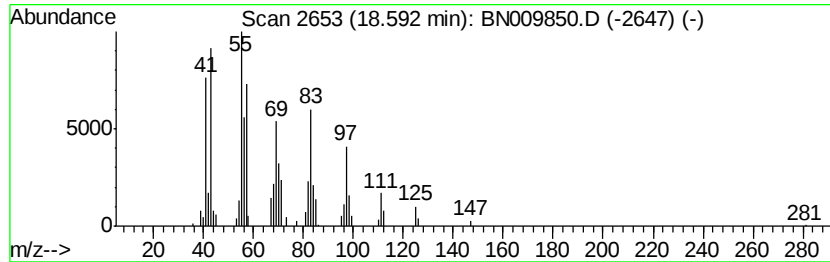
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 (DEL) Alkane: Straight-Chai... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.59	2.35 ng/ul	193342	Phenanthrene-d10	16.79

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Nonadecene	266	C19H38	018435-45-5	91
2			Cyclohexadecane	224	C16H32	000295-65-8	91
3			n-Heptadecanol-1	256	C17H36O	001454-85-9	91
4			E-15-Heptadecenal	252	C17H32O	1000130-97-9	91
5			E-14-Hexadecenal	238	C16H30O	330207-53-9	91



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Data File : BN009850.D
Acq On : 22 Feb 2020 17:45
Operator : CG/JU
Sample : L1516-12
Misc :
ALS Vial : 39 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
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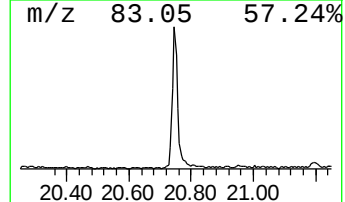
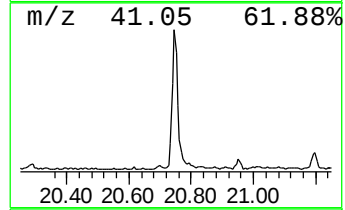
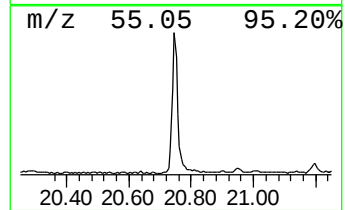
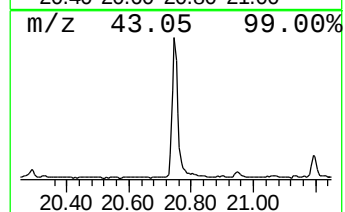
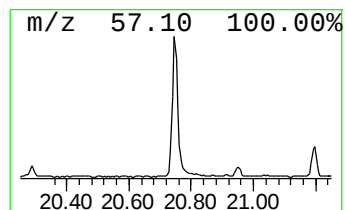
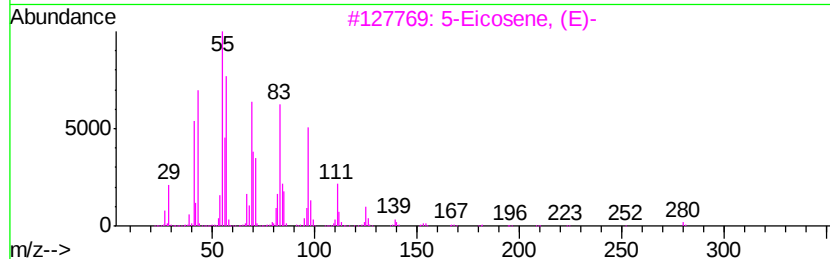
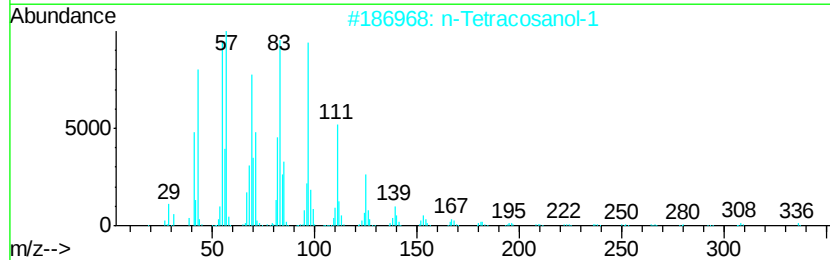
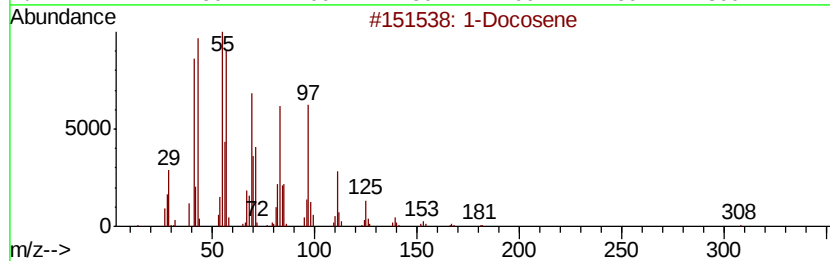
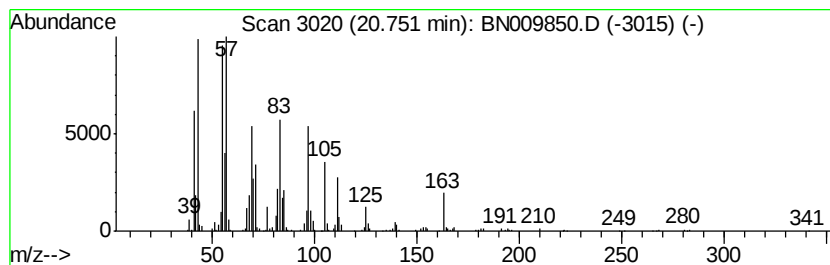
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 (DEL) Alkane: Straight-Chai... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.75	9.61 ng/ul	863419	Chrysene-d12	21.00

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Docosene	308	C22H44	001599-67-3	95
2			n-Tetracosanol-1	354	C24H50O	000506-51-4	91
3			5-Eicosene, (E)-	280	C20H40	074685-30-6	91
4			1-Heneicosanol	312	C21H44O	015594-90-8	91
5			Trifluoroacetic acid, pentadecyl...	324	C17H31F3O2	959010-23-2	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN022220\
Data File : BN009850.D
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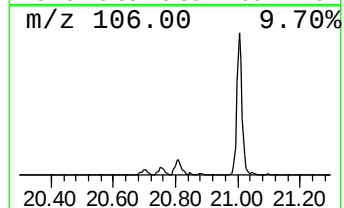
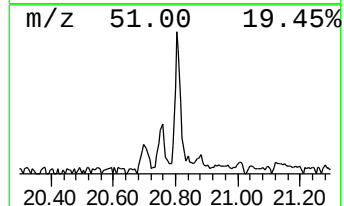
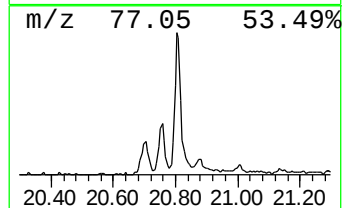
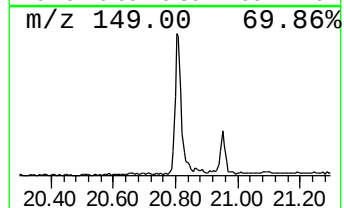
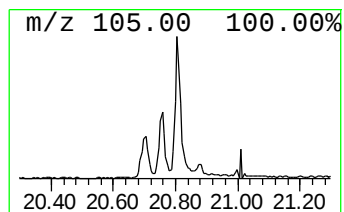
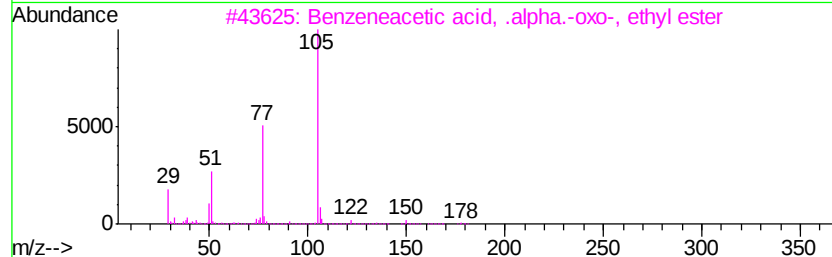
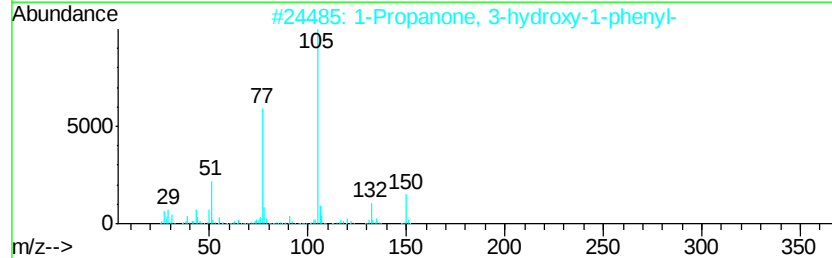
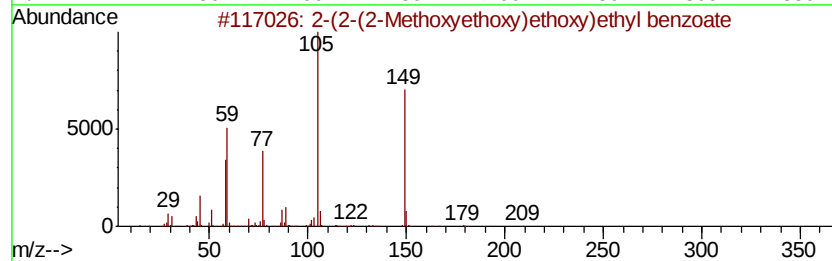
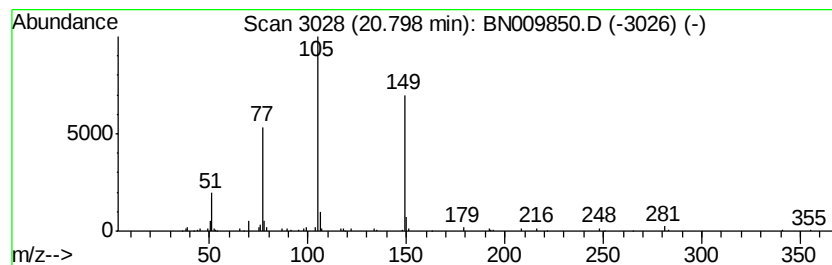
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 2-(2-(2-Methoxyethoxy)ethox... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.80	2.59 ng/ul	232833	Chrysene-d12	21.00

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		2-(2-(2-Methoxyethoxy)ethoxy)eth...	268	C14H20O5	1000366-99-1	72
2		1-Propanone, 3-hydroxy-1-phenyl-	150	C9H10O2	005650-41-9	49
3		Benzeneacetic acid, .alpha.-oxo-...	178	C10H10O3	001603-79-8	47
4		Benzoic acid, ethyl ester	150	C9H10O2	000093-89-0	47
5		Benzoic acid, ethyl ester	150	C9H10O2	000093-89-0	47



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN022220\
Data File : BN009850.D
Acq On : 22 Feb 2020 17:45
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ClientSampleId :
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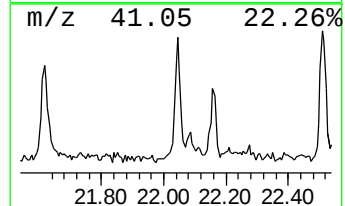
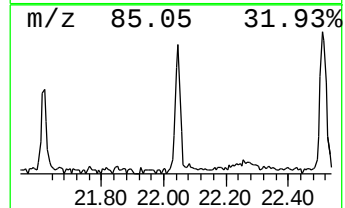
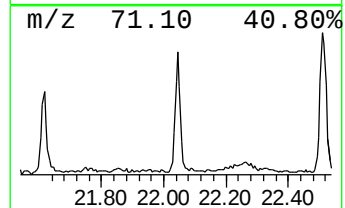
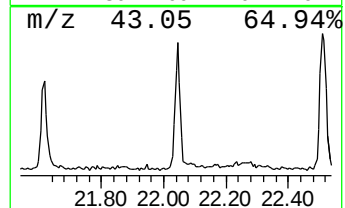
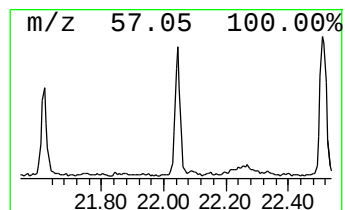
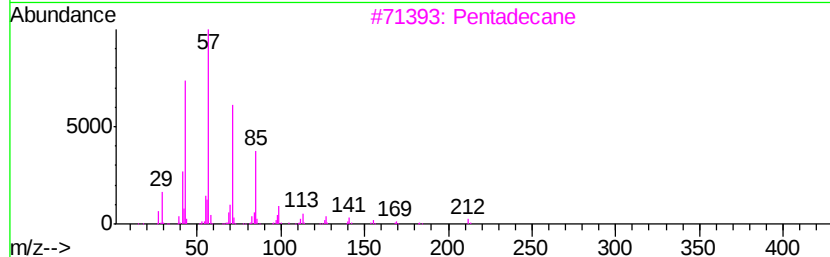
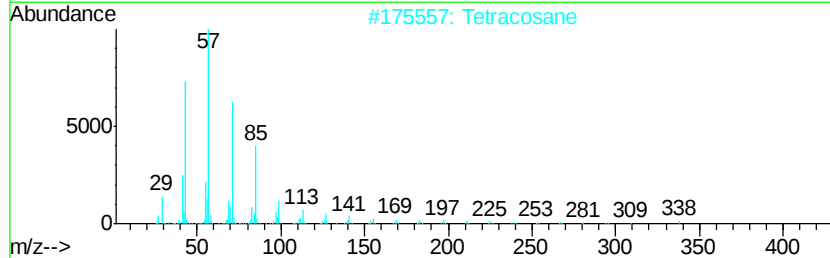
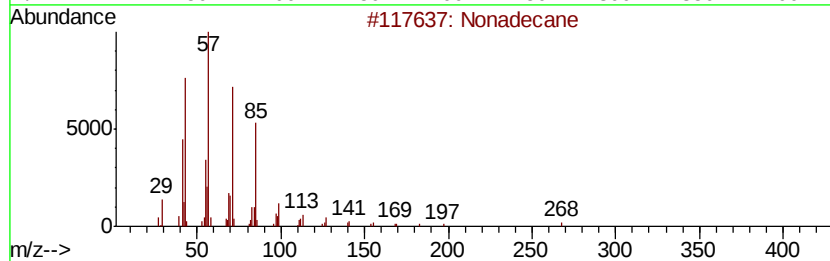
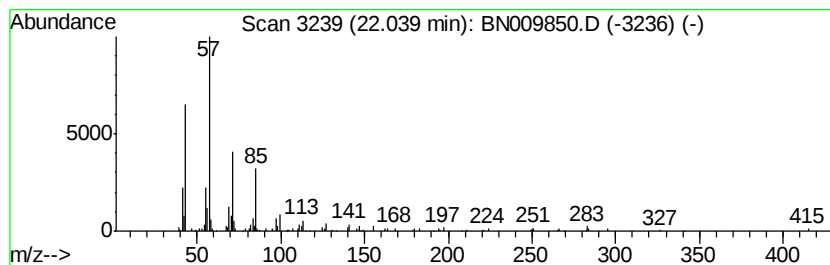
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 (DEL) Alkane: Straight-Chai... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.04	2.06 ng/ul	184742	Chrysene-d12	21.00

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Nonadecane	268	C19H40	000629-92-5	93
2		Tetracosane	338	C24H50	000646-31-1	87
3		Pentadecane	212	C15H32	000629-62-9	83
4		Eicosane	282	C20H42	000112-95-8	83
5		Nonadecane	268	C19H40	000629-92-5	83



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN022220\
Data File : BN009850.D
Acq On : 22 Feb 2020 17:45
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Misc :
ALS Vial : 39 Sample Multiplier: 1

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ClientSampleId :
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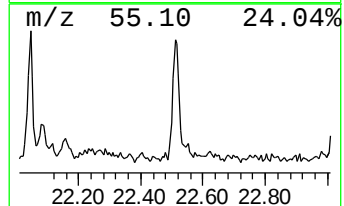
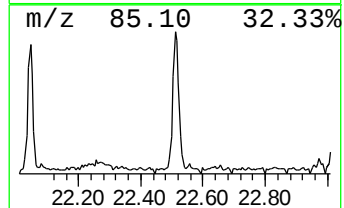
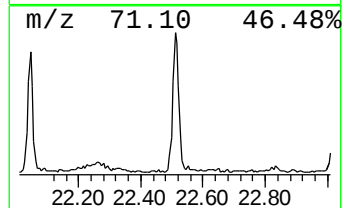
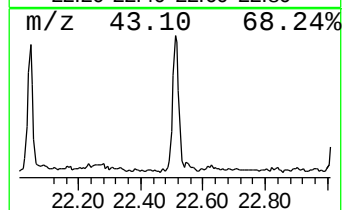
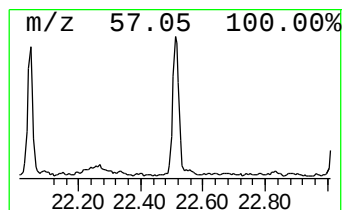
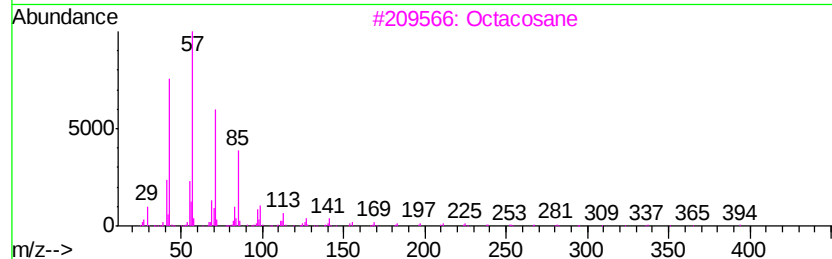
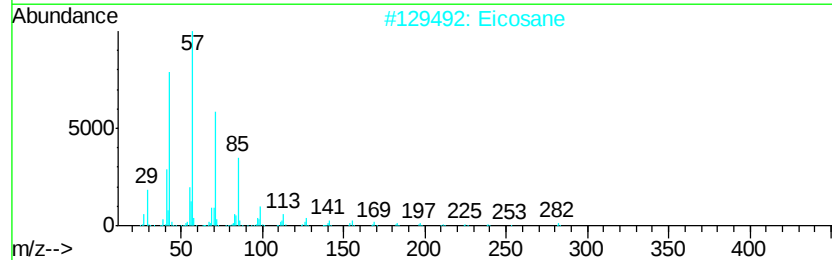
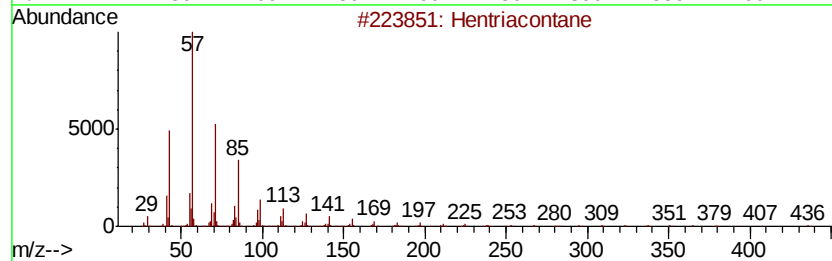
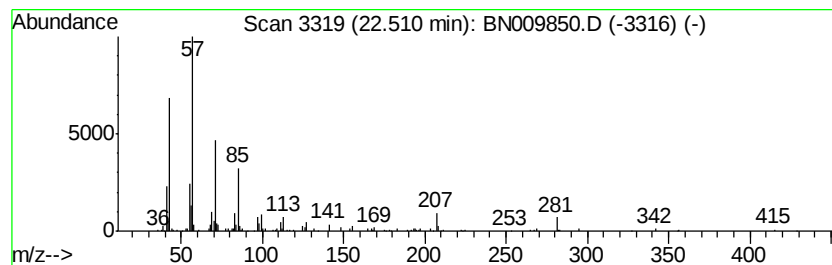
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 (DEL) Alkane: Straight-Chai... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.51	2.45 ng/ul	218465	Perylene-d12	23.10

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hentriacontane	437	C31H64	000630-04-6	68
2			Eicosane	282	C20H42	000112-95-8	68
3			Octacosane	394	C28H58	000630-02-4	64
4			Hexacosane	366	C26H54	000630-01-3	64
5			Docosane	310	C22H46	000629-97-0	64



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN022220\
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Acq On : 22 Feb 2020 17:45
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Misc :
ALS Vial : 39 Sample Multiplier: 1

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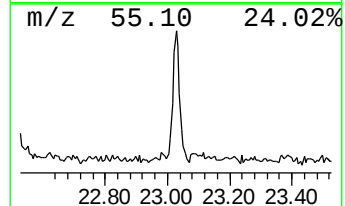
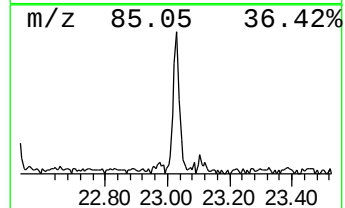
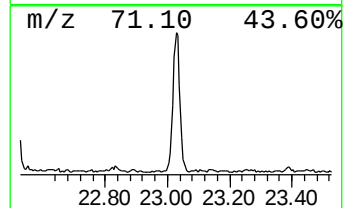
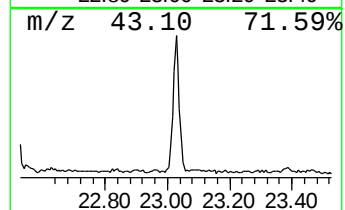
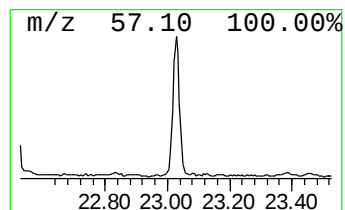
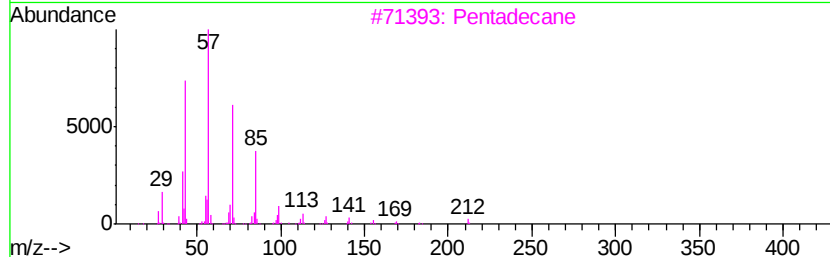
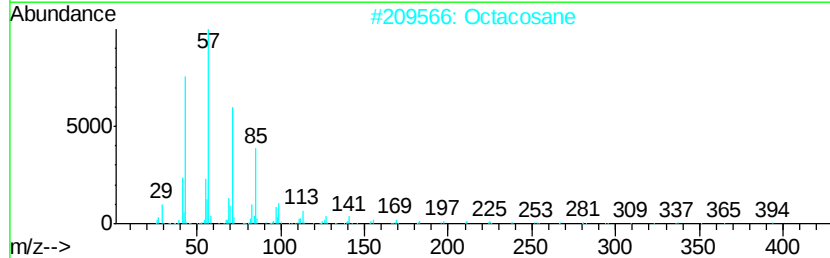
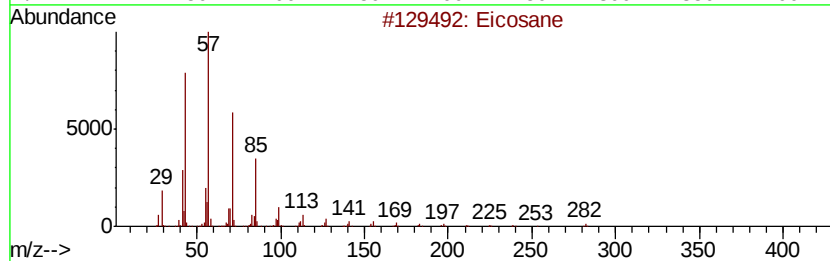
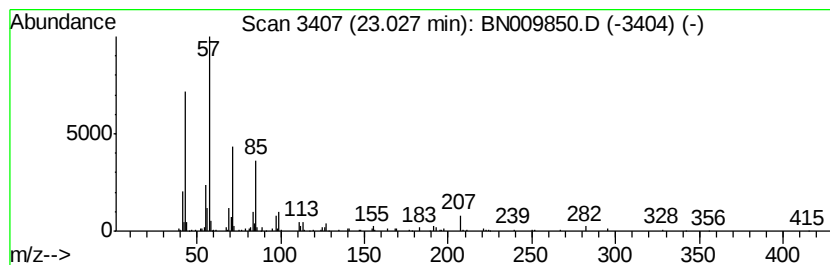
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 (DEL) Alkane: Straight-Chai... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.03	3.29 ng/ul	293302	Perylene-d12	23.10

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Eicosane	282	C20H42	000112-95-8	98
2		Octacosane	394	C28H58	000630-02-4	87
3		Pentadecane	212	C15H32	000629-62-9	80
4		Hexadecane	226	C16H34	000544-76-3	80
5		Sulfurous acid, butyl undecyl ester	292	C15H32O3S	1000309-17-8	80



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN022220\
Data File : BN009850.D
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ALS Vial : 39 Sample Multiplier: 1

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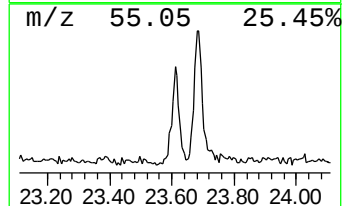
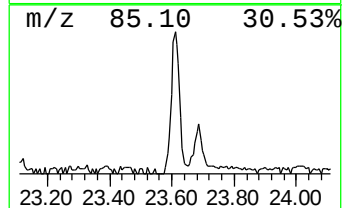
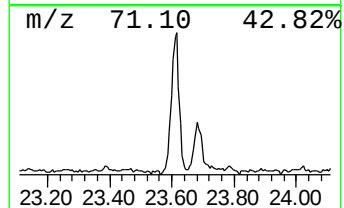
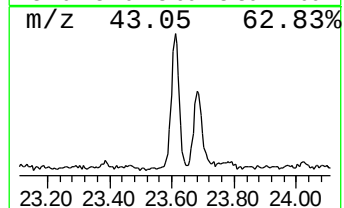
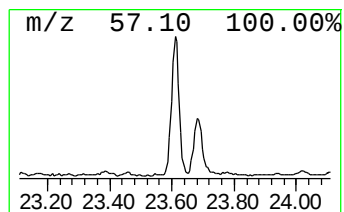
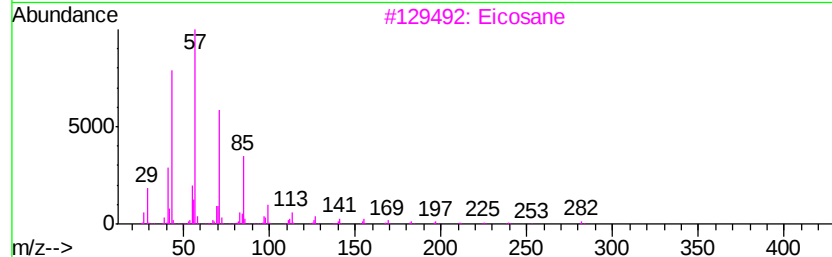
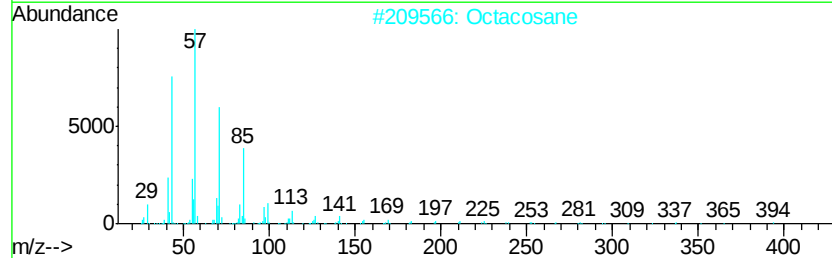
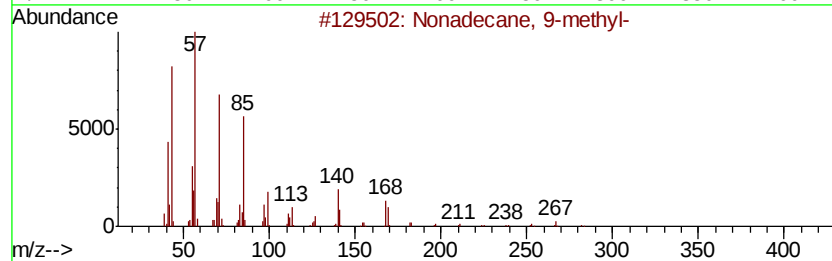
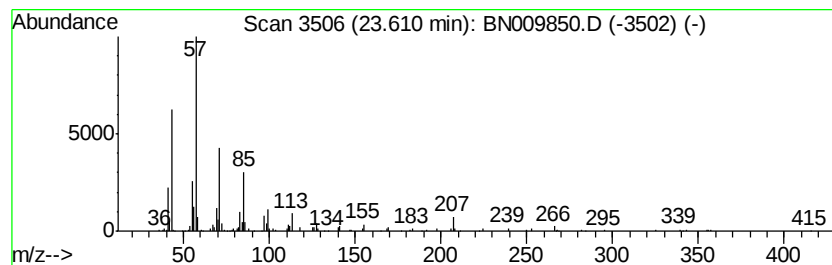
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 9 (DEL) Alkane: Straight-Chai... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.61	3.06 ng/ul	272839	Perylene-d12	23.10

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Nonadecane, 9-methyl-	282	C20H42	013287-24-6	89
2		Octacosane	394	C28H58	000630-02-4	87
3		Eicosane	282	C20H42	000112-95-8	87
4		Octadecane, 3-methyl-	268	C19H40	006561-44-0	83
5		Heptacosane	380	C27H56	000593-49-7	81



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN022220\
Data File : BN009850.D
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Sample : L1516-12
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ALS Vial : 39 Sample Multiplier: 1

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ClientSampleId :
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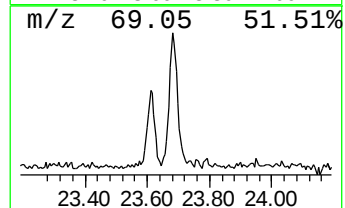
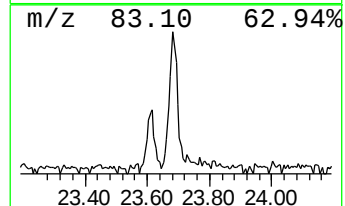
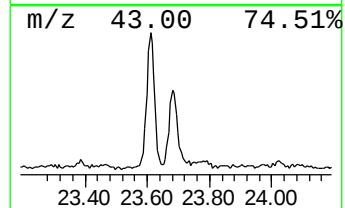
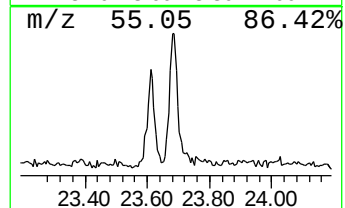
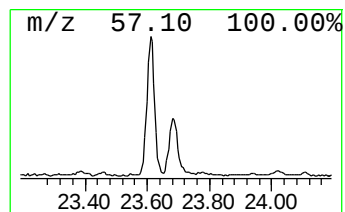
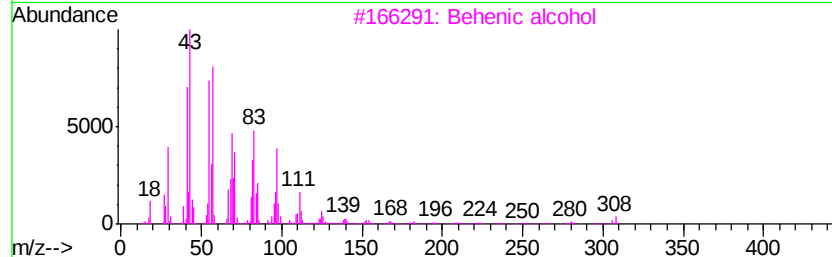
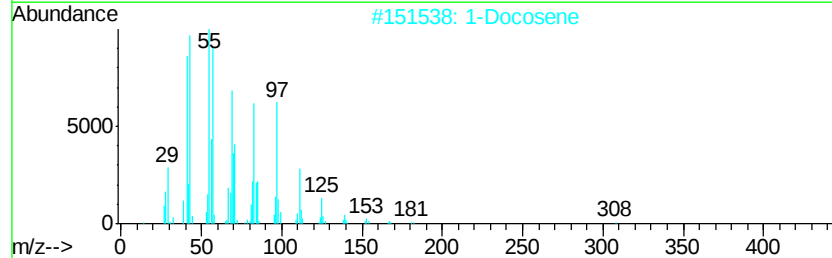
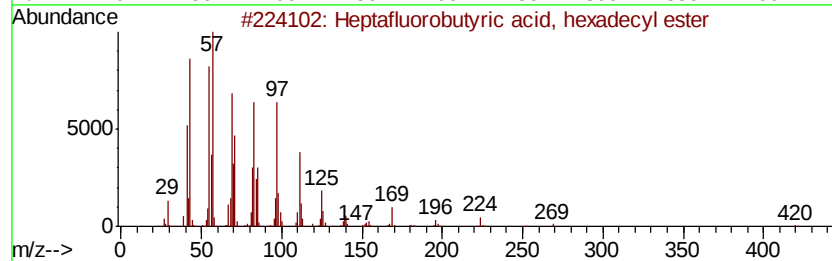
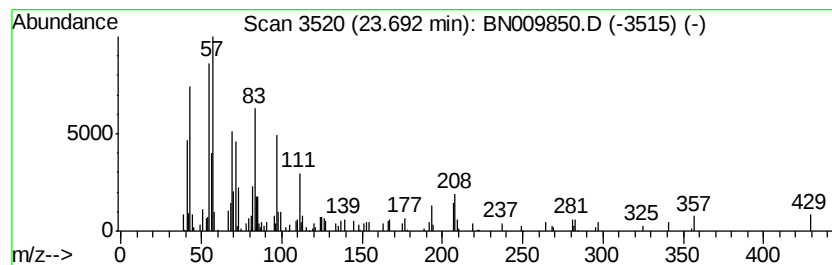
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 Heptafluorobutyric acid, he... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.69	2.43 ng/ul	216302	Perylene-d12	23.10

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heptafluorobutyric acid, hexadec...	438	C20H33F7O2	006385-15-5	81
2		1-Docosene	308	C22H44	001599-67-3	81
3		Behenic alcohol	326	C22H46O	000661-19-8	74
4		Trichloroacetic acid, pentadecyl...	372	C17H31Cl3O2	074339-53-0	74
5		Pentafluoropropionic acid, hexad...	388	C19H33F5O2	006222-07-7	74



Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN022220\
Data File : BN009850.D
Acq On : 22 Feb 2020 17:45
Operator : CG/JU
Sample : L1516-12
Misc :
ALS Vial : 39 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
DBD30

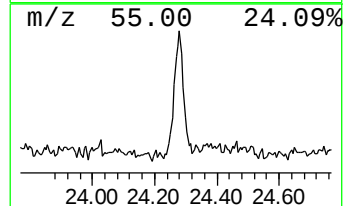
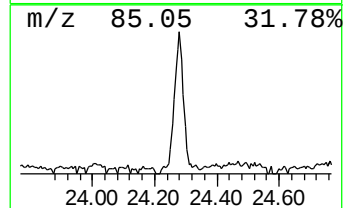
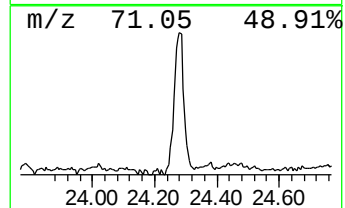
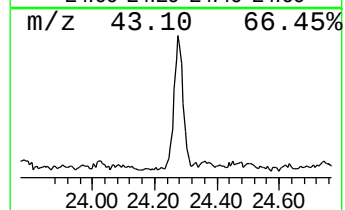
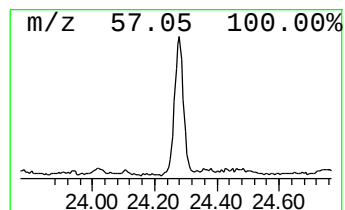
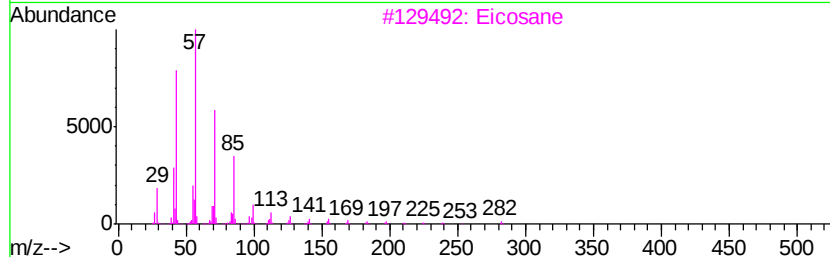
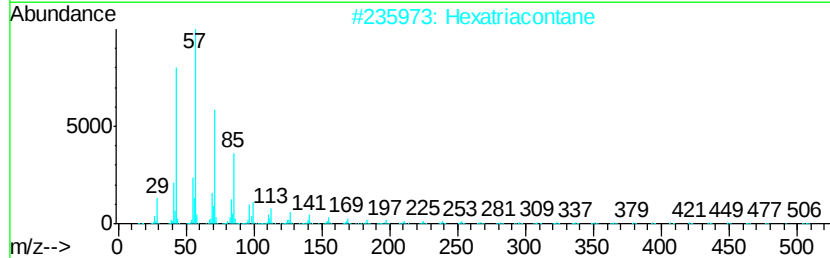
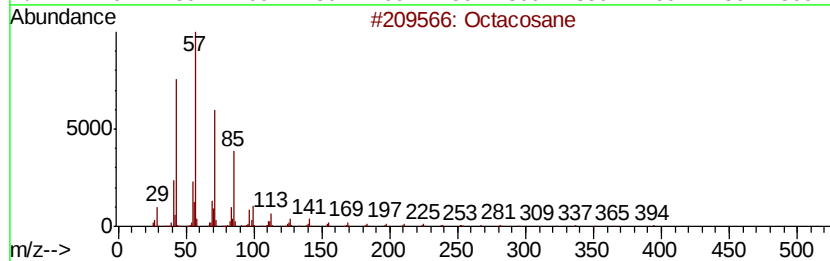
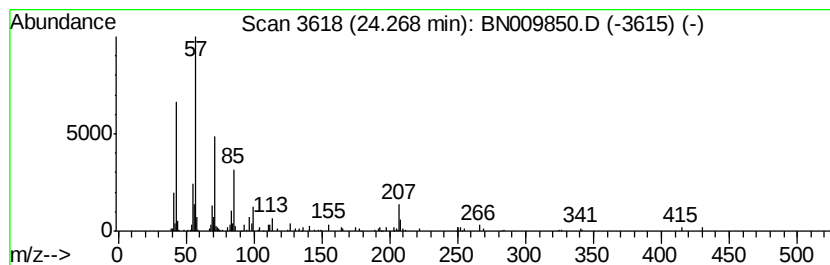
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Quant Title : SVOA CALIBRATION

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

Peak Number 11 (DEL) Alkane: Straight-Chai... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.27	2.81 ng/ul	250397	Perylene-d12	23.10

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octacosane	394	C28H58	000630-02-4	76
2			Hexatriacontane	507	C36H74	000630-06-8	68
3			Eicosane	282	C20H42	000112-95-8	68
4			Tetratetracontane	619	C44H90	007098-22-8	64
5			Hexacosane	366	C26H54	000630-01-3	64



Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN022220\
Data File : BN009850.D
Acq On : 22 Feb 2020 17:45
Operator : CG/JU
Sample : L1516-12
Misc :
ALS Vial : 39 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
DBD30

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN021220MA.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
unknown-01	12.95	6.2	ng/ul	436269	3	14.03	1405120	20.0
n-Hexadecanoic acid	17.72	5.1	ng/ul	422330	4	16.79	1647230	20.0
(DEL) Alkane: Str...	18.59	2.4	ng/ul	193342	4	16.79	1647230	20.0
(DEL) Alkane: Str...	20.75	9.6	ng/ul	863419	5	21.00	1796670	20.0
2-(2-(2-Methoxyet...	20.80	2.6	ng/ul	232833	5	21.00	1796670	20.0
(DEL) Alkane: Str...	22.04	2.1	ng/ul	184742	5	21.00	1796670	20.0
(DEL) Alkane: Str...	22.51	2.5	ng/ul	218465	6	23.10	1782340	20.0
(DEL) Alkane: Str...	23.03	3.3	ng/ul	293302	6	23.10	1782340	20.0
(DEL) Alkane: Str...	23.61	3.1	ng/ul	272839	6	23.10	1782340	20.0
Heptafluorobutyri...	23.69	2.4	ng/ul	216302	6	23.10	1782340	20.0
(DEL) Alkane: Str...	24.27	2.8	ng/ul	250397	6	23.10	1782340	20.0