

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN030420\
 Data File : BN010097.D
 Acq On : 05 Mar 2020 02:27
 Operator : CG/JU
 Sample : L1617-11
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 DBDS5

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	3.193	34	35	39	rVB3	3422	3428	0.13%	0.010%
2	3.687	115	119	126	rVB6	5532	11610	0.45%	0.036%
3	4.058	179	182	184	rVB3	2630	2891	0.11%	0.009%
4	4.293	219	222	224	rVB4	2749	2731	0.11%	0.008%
5	4.569	265	269	275	rBV5	5888	9015	0.35%	0.028%
6	5.146	364	367	370	rBV3	2490	3529	0.14%	0.011%
7	5.540	430	434	437	rVB5	2513	4359	0.17%	0.013%
8	5.764	467	472	473	rBV3	2047	3177	0.12%	0.010%
9	6.563	602	608	622	rBV	555779	1004324	38.72%	3.073%
10	6.716	629	634	651	rVV	392713	643889	24.82%	1.970%
11	6.899	659	665	677	rVV	558059	916407	35.33%	2.804%
12	7.352	736	742	752	rBV	352761	580797	22.39%	1.777%
13	7.452	754	759	761	rBV3	6267	8496	0.33%	0.026%
14	8.075	857	865	878	rBV	631164	1103200	42.53%	3.376%
15	8.216	886	889	892	rVB3	2421	3355	0.13%	0.010%
16	8.499	931	937	951	rBV	537672	950804	36.65%	2.910%
17	8.675	964	967	968	rBV3	3634	3251	0.13%	0.010%
18	8.934	1003	1011	1016	rBV2	12968	16994	0.66%	0.052%
19	9.210	1051	1058	1074	rBV	476123	824377	31.78%	2.523%
20	9.310	1074	1075	1080	rVB5	2331	2814	0.11%	0.009%
21	9.746	1140	1149	1168	rBV	584735	1125546	43.39%	3.444%
22	10.099	1202	1209	1222	rBV	514519	849372	32.74%	2.599%
23	10.263	1230	1237	1257	rBV	483439	1086324	41.88%	3.324%
24	11.051	1368	1371	1377	rVB3	1696	2798	0.11%	0.009%
25	11.716	1479	1484	1489	rBV3	4888	9590	0.37%	0.029%
26	12.334	1587	1589	1594	rBV3	2266	3829	0.15%	0.012%
27	12.446	1603	1608	1611	rBV2	2169	3189	0.12%	0.010%
28	12.840	1669	1675	1681	rBV4	4517	7851	0.30%	0.024%
29	12.910	1683	1687	1693	rVB2	9884	14192	0.55%	0.043%
30	13.428	1769	1775	1780	rBV	1104954	1570459	60.54%	4.806%
31	13.475	1780	1783	1791	rVB	110914	171305	6.60%	0.524%
32	13.687	1813	1819	1834	rBV	1272817	1872302	72.18%	5.730%
33	13.910	1853	1857	1859	rBV3	2701	3356	0.13%	0.010%
34	13.934	1859	1861	1863	rVV2	3096	3416	0.13%	0.010%

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Integration Parameters: LSCINT.P

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 Title : SVOA CALIBRATION

35	13.998	1866	1872	1886	rVV2	758452	1118297	43.11%	3.422%
36	14.098	1886	1889	1892	rVB3	3674	3950	0.15%	0.012%
37	14.263	1911	1917	1936	rBV	343572	785836	30.29%	2.405%
38	14.463	1949	1951	1956	rVB5	2867	3795	0.15%	0.012%
39	14.898	2021	2025	2030	rBV4	7547	10990	0.42%	0.034%
40	15.010	2037	2044	2053	rBV	1504539	2268488	87.45%	6.942%
41	15.069	2053	2054	2059	rVB3	2629	2857	0.11%	0.009%
42	15.175	2066	2072	2082	rBV	502629	806117	31.08%	2.467%
43	15.251	2082	2085	2089	rVB3	20110	28542	1.10%	0.087%
44	15.334	2097	2099	2102	rBV	1811	2610	0.10%	0.008%
45	15.763	2169	2172	2176	rBV5	3987	6608	0.25%	0.020%
46	15.798	2176	2178	2181	rVB2	5082	4363	0.17%	0.013%
47	15.934	2196	2201	2203	rBV4	2495	3140	0.12%	0.010%
48	16.010	2211	2214	2218	rVB2	5645	6569	0.25%	0.020%
49	16.169	2239	2241	2244	rBV4	3632	3727	0.14%	0.011%
50	16.216	2244	2249	2251	rBV4	2933	3982	0.15%	0.012%
51	16.516	2297	2300	2303	rBV4	3089	3476	0.13%	0.011%
52	16.563	2303	2308	2312	rBV4	6290	11249	0.43%	0.034%
53	16.763	2336	2342	2352	rVB	878967	1242147	47.88%	3.801%
54	16.863	2352	2359	2371	rBV	1717852	2399654	92.51%	7.344%
55	17.004	2378	2383	2387	rVB5	6011	11148	0.43%	0.034%
56	17.181	2405	2413	2419	rBV2	34225	56916	2.19%	0.174%
57	17.428	2451	2455	2456	rBV3	3313	3755	0.14%	0.011%
58	17.692	2495	2500	2509	rBV	153228	259429	10.00%	0.794%
59	17.969	2544	2547	2549	rBV3	7303	8083	0.31%	0.025%
60	17.992	2549	2551	2553	rVB3	3183	2823	0.11%	0.009%
61	18.128	2569	2574	2578	rVB3	14691	19078	0.74%	0.058%
62	18.386	2616	2618	2621	rVV3	2798	2700	0.10%	0.008%
63	18.422	2621	2624	2626	rBV3	3224	4336	0.17%	0.013%
64	18.498	2634	2637	2641	rBV5	5032	6354	0.24%	0.019%
65	18.563	2645	2648	2657	rVB7	9300	18780	0.72%	0.057%
66	18.633	2657	2660	2662	rBV2	6026	6762	0.26%	0.021%
67	18.686	2665	2669	2673	rBV3	5819	9268	0.36%	0.028%
68	18.804	2686	2689	2693	rBV	14042	19824	0.76%	0.061%
69	18.839	2693	2695	2697	rBV3	4579	3994	0.15%	0.012%
70	18.992	2717	2721	2727	rBV3	32831	64529	2.49%	0.197%
71	19.063	2730	2733	2741	rVB2	17733	29413	1.13%	0.090%

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 Title : SVOA CALIBRATION

72	19.175	2746	2752	2764	rVV	1923995	2594073	100.00%	7.938%
73	19.357	2781	2783	2785	rBV3	4260	4329	0.17%	0.013%
74	19.498	2805	2807	2810	rBV4	5201	6331	0.24%	0.019%
75	19.528	2810	2812	2814	rBV3	5561	4702	0.18%	0.014%
76	19.733	2842	2847	2850	rBV7	19608	32430	1.25%	0.099%
77	20.145	2914	2917	2921	rBV6	13536	17289	0.67%	0.053%
78	20.680	3003	3008	3011	rBV	70058	118510	4.57%	0.363%
79	20.722	3011	3015	3022	rVV2	525442	734560	28.32%	2.248%
80	20.786	3022	3026	3033	rVV	281393	386076	14.88%	1.181%
81	20.986	3055	3060	3067	rBV2	998004	1297384	50.01%	3.970%
82	21.110	3078	3081	3086	rBV2	77649	103728	4.00%	0.317%
83	21.169	3088	3091	3095	rBV3	54654	66019	2.54%	0.202%
84	21.586	3159	3162	3168	rBV2	81202	130845	5.04%	0.400%
85	22.016	3232	3235	3239	rVB2	147388	162664	6.27%	0.498%
86	22.480	3310	3314	3318	rBV	202141	257233	9.92%	0.787%
87	22.951	3388	3394	3399	rBV	1415261	2411727	92.97%	7.380%
88	22.992	3399	3401	3407	rVB	240383	302168	11.65%	0.925%
89	23.080	3410	3416	3424	rVB	694541	1274336	49.12%	3.900%
90	23.568	3495	3499	3506	rVB	219350	350904	13.53%	1.074%
91	24.233	3607	3612	3620	rVB	189021	355251	13.69%	1.087%

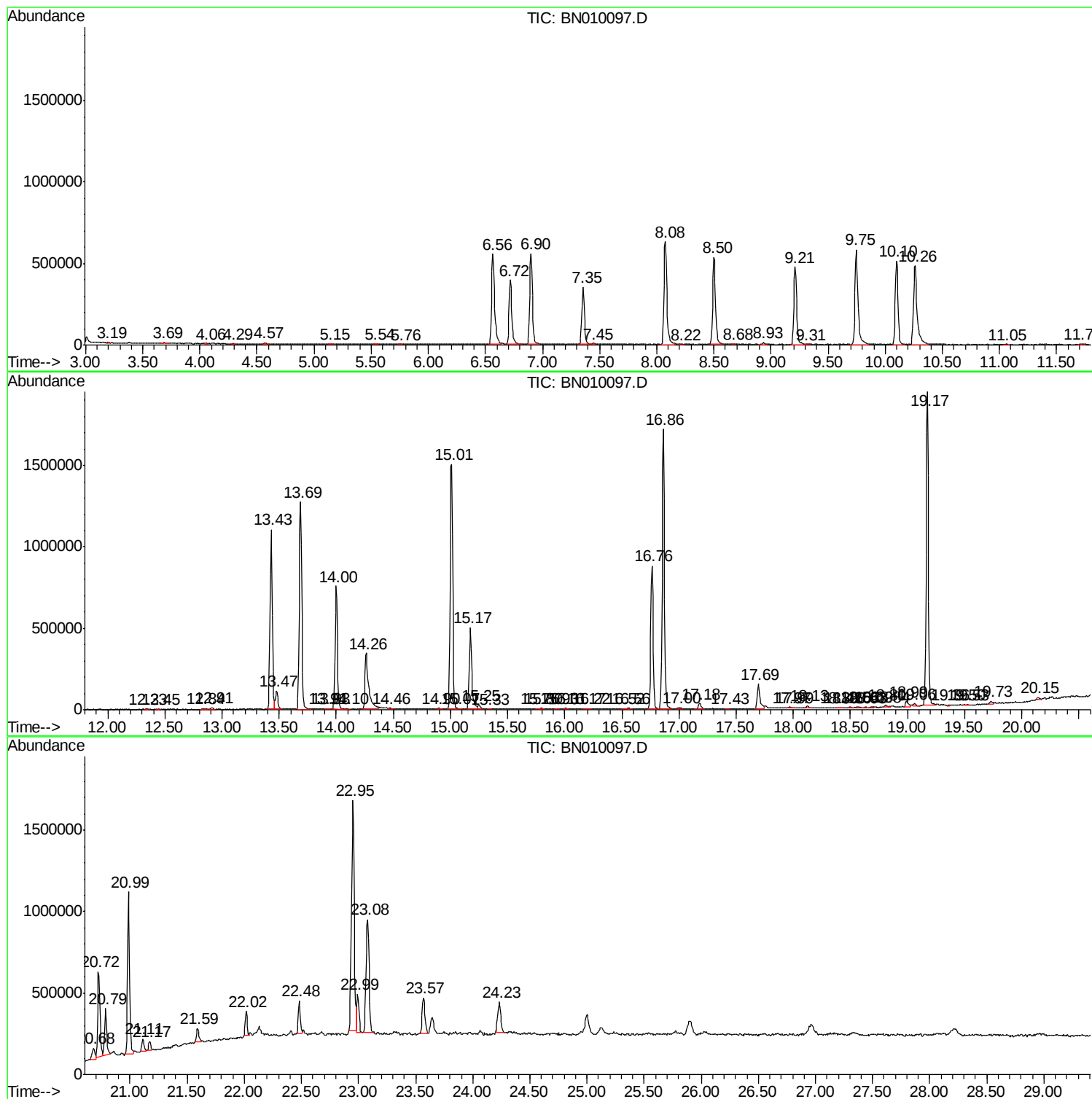
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN021220MA.M
Quant Title : SVOA CALIBRATION

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



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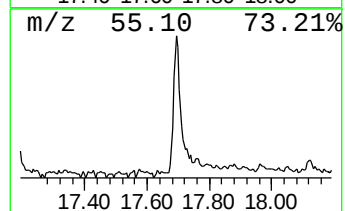
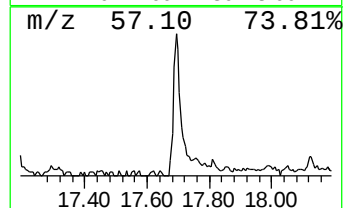
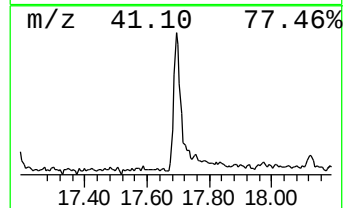
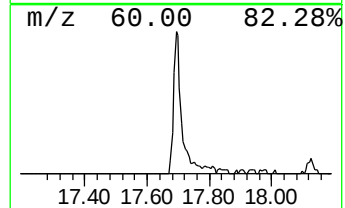
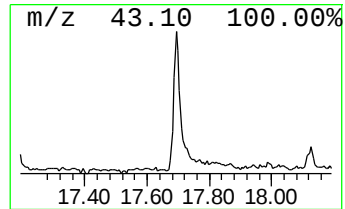
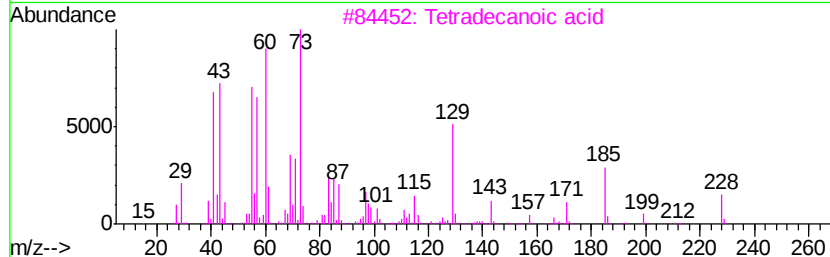
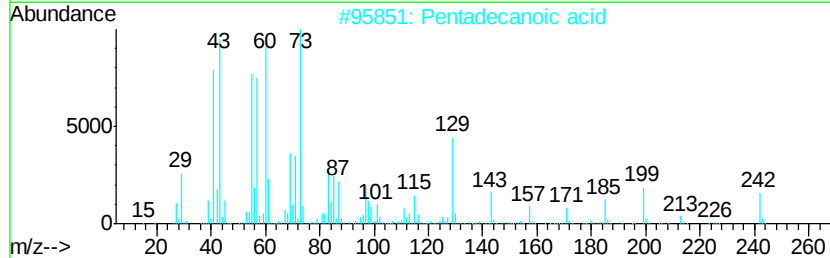
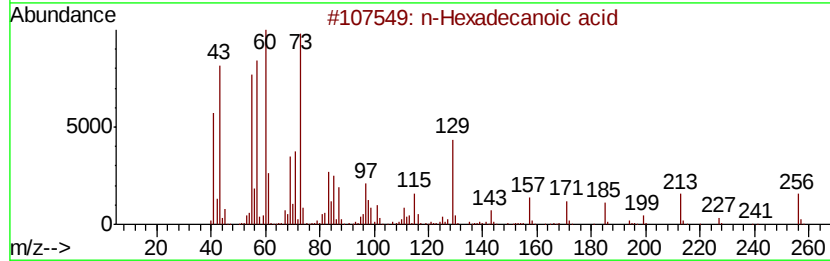
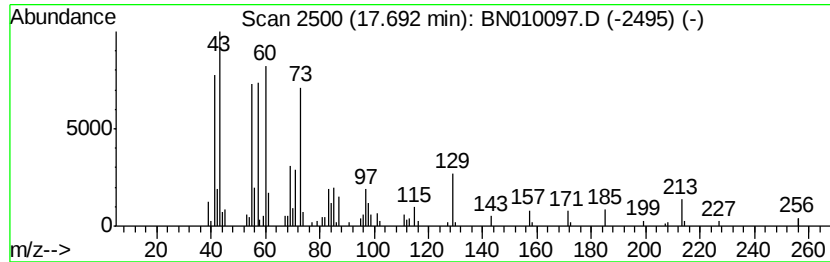
Instrument :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN021220MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 n-Hexadecanoic acid Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.	
17.69	4.18 ng/ul	259429	Phenanthrene-d10	16.76	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
2	Pentadecanoic acid	242	C15H30O2	001002-84-2	91
3	Tetradecanoic acid	228	C14H28O2	000544-63-8	90
4	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	78
5	Tridecanoic acid	214	C13H26O2	000638-53-9	72



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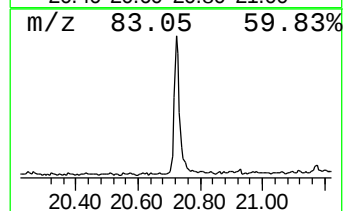
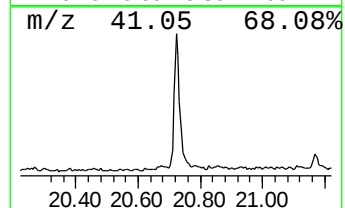
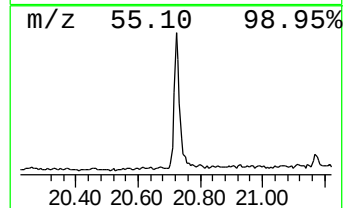
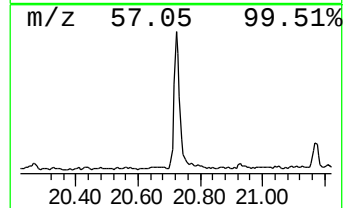
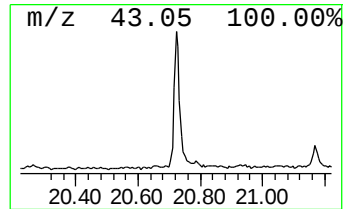
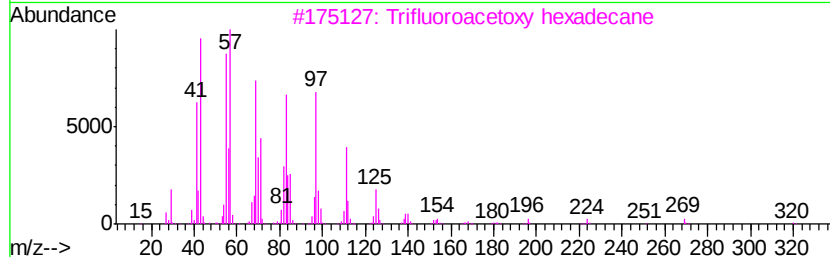
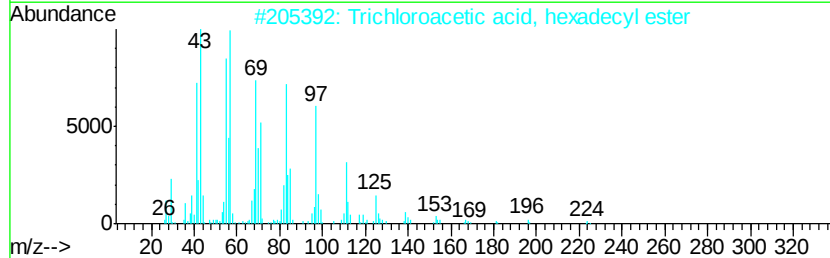
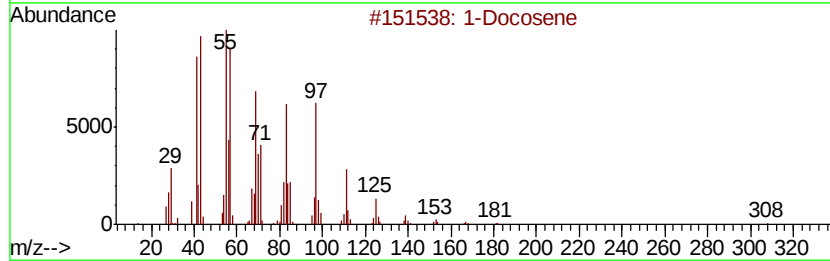
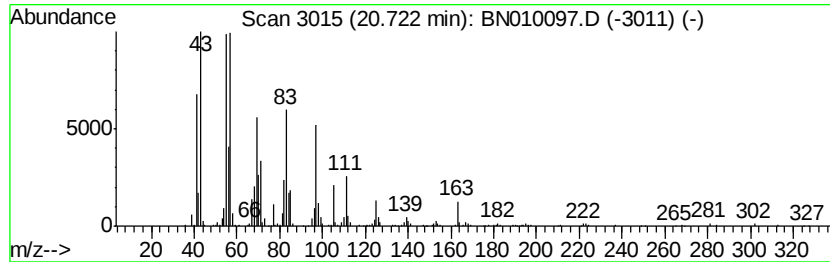
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN021220MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 (DEL) Alkane: Straight-Chai... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.72	11.32 ng/ul	734560	Chrysene-d12	20.99

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1-Docosene	308	C22H44	001599-67-3	91
2		Trichloroacetic acid, hexadecyl ...	386	C18H33Cl3O2	074339-54-1	87
3		Trifluoroacetoxv hexadecane	338	C18H33F3O2	006222-03-3	87
4		1-Heneicosanol	312	C21H44O	015594-90-8	87
5		n-Tetracosanol-1	354	C24H50O	000506-51-4	87



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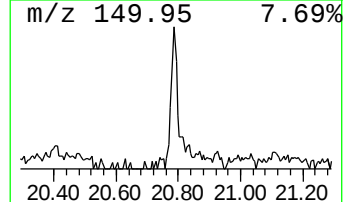
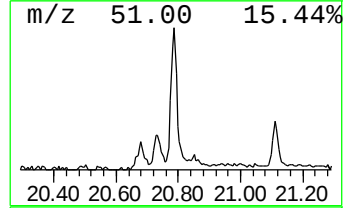
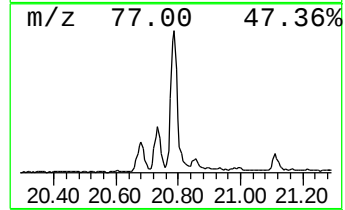
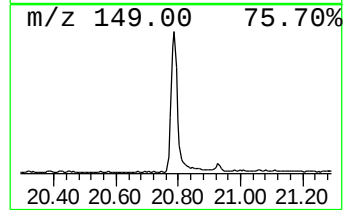
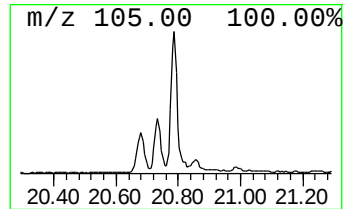
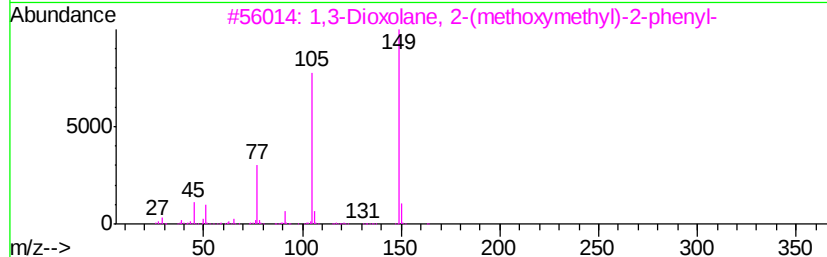
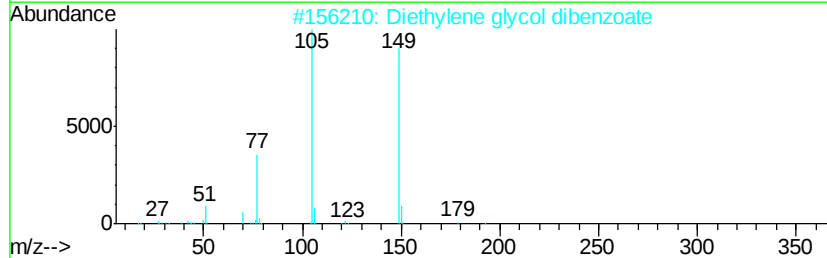
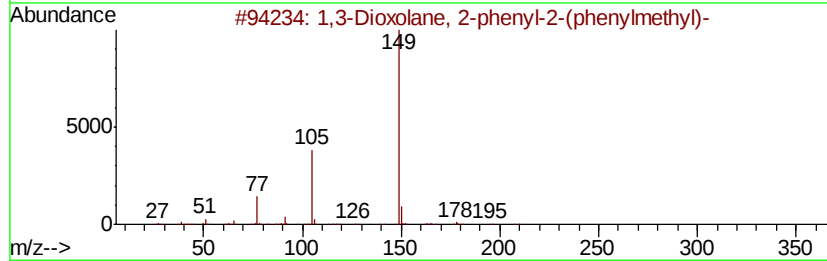
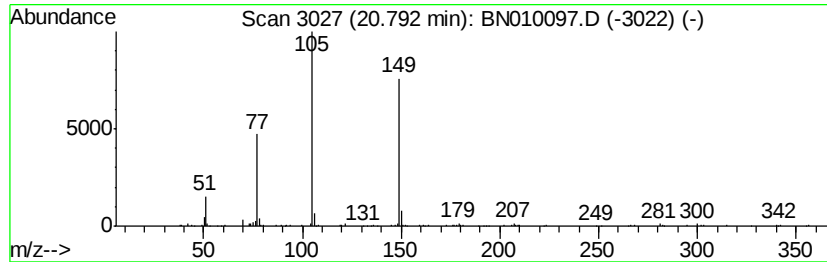
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN021220MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 1,3-Dioxolane, 2-phenyl-2-(... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.79	5.95 ng/ul	386076	Chrysene-d12	20.99

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1,3-Dioxolane, 2-phenyl-2-(phenyl...	240	C16H16O2	004362-19-0	78
2		Diethylene glycol dibenzoate	314	C18H18O5	000120-55-8	78
3		1,3-Dioxolane, 2-(methoxymethyl)...	194	C11H14O3	1000156-71-2	72
4		1-Propanone, 3-hydroxy-1-phenyl-	150	C9H10O2	005650-41-9	47
5		Benzeneacetic acid, .alpha.-oxo-...	178	C10H10O3	001603-79-8	47



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ClientSampleId :
DBDS5

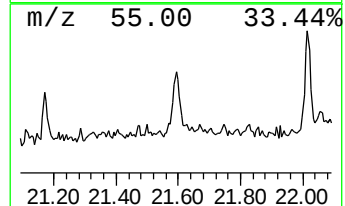
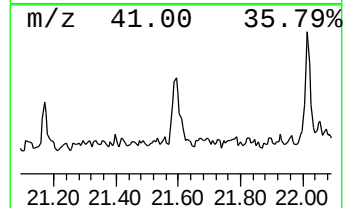
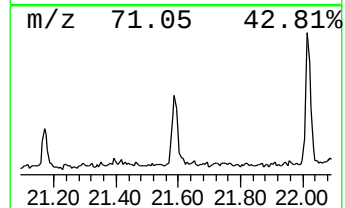
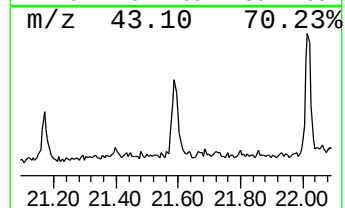
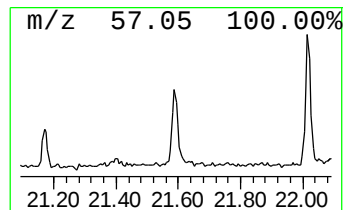
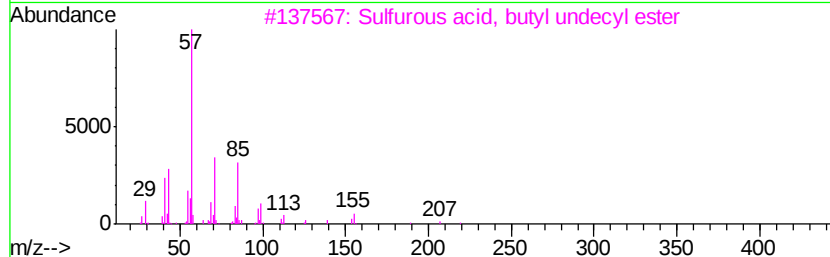
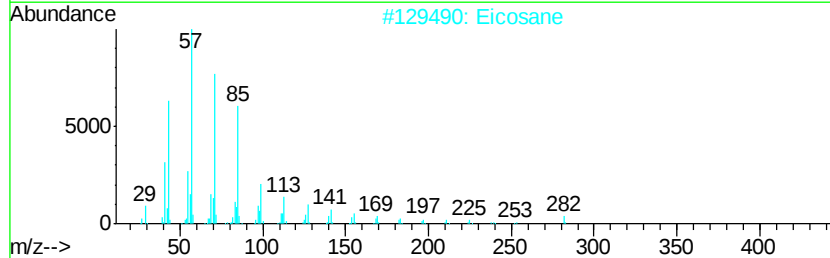
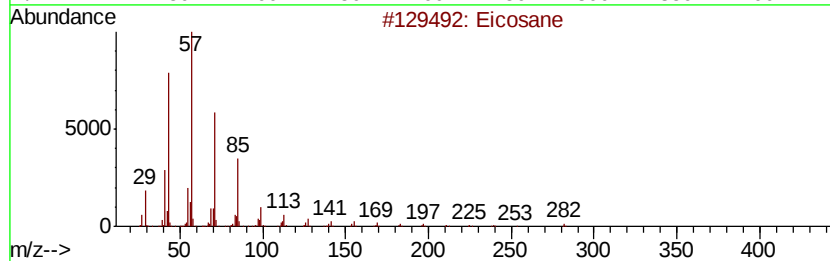
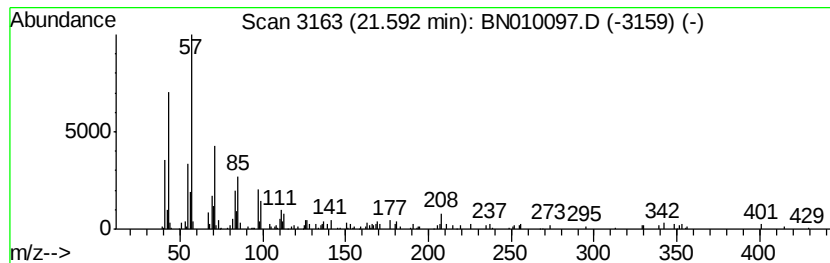
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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.59	2.02 ng/ul	130845	Chrysene-d12	20.99

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Eicosane	282	C20H42	000112-95-8	94
2		Eicosane	282	C20H42	000112-95-8	93
3		Sulfurous acid, butyl undecyl ester	292	C15H32O3S	1000309-17-8	83
4		Octadecane, 2-methyl-	268	C19H40	001560-88-9	80
5		Heneicosane, 11-decyl-	437	C31H64	055320-06-4	80



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN030420\
Data File : BN010097.D
Acq On : 05 Mar 2020 02:27
Operator : CG/JU
Sample : L1617-11
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
DBDS5

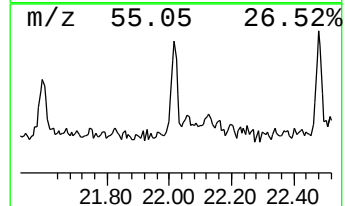
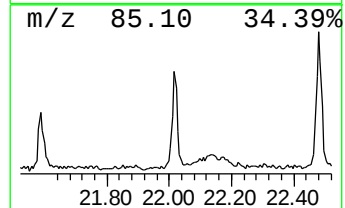
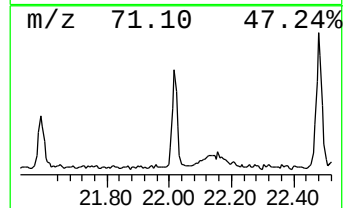
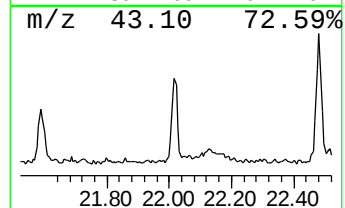
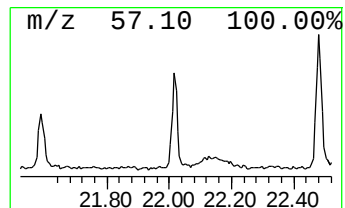
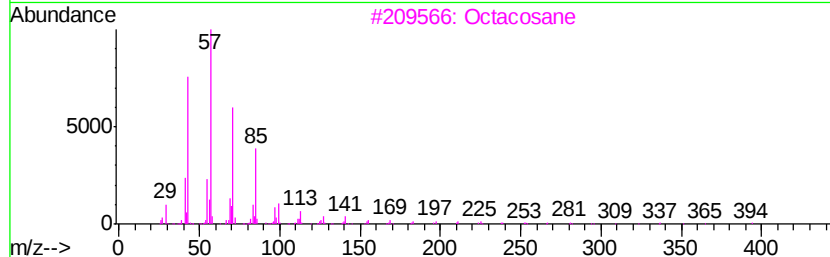
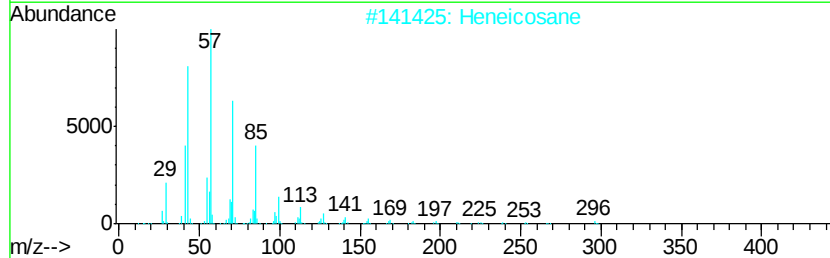
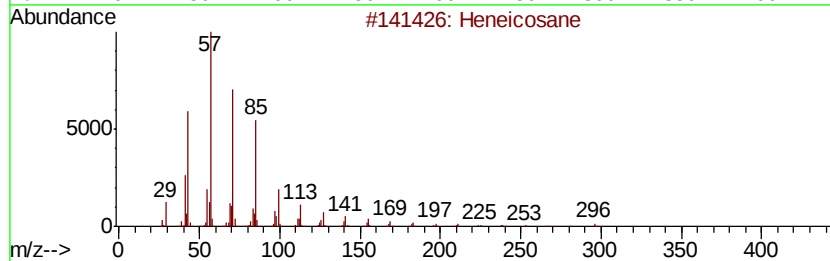
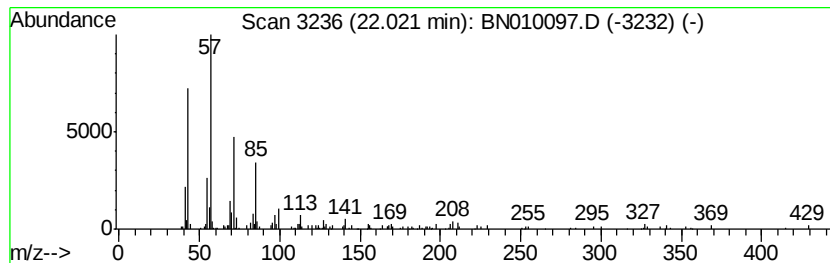
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 (DEL) Alkane: Straight-Chai... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.02	2.51 ng/ul	162664	Chrysene-d12	20.99

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heneicosane		296	C21H44	000629-94-7	94
2	Heneicosane		296	C21H44	000629-94-7	93
3	Octacosane		394	C28H58	000630-02-4	81
4	Tetracosane		338	C24H50	000646-31-1	81
5	Heneicosane		296	C21H44	000629-94-7	76



Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN030420\
Data File : BN010097.D
Acq On : 05 Mar 2020 02:27
Operator : CG/JU
Sample : L1617-11
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
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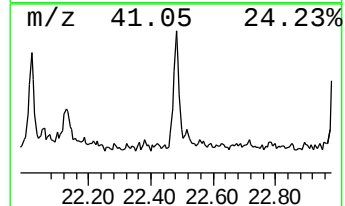
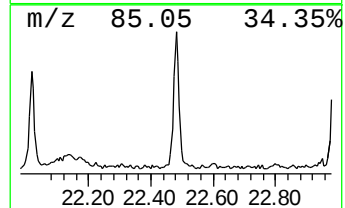
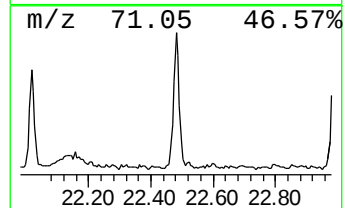
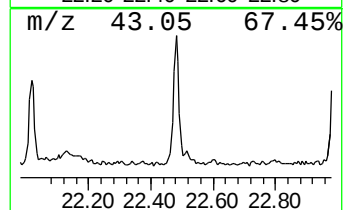
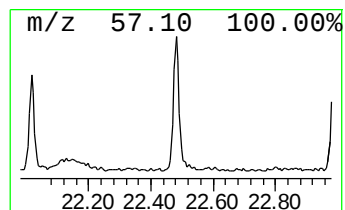
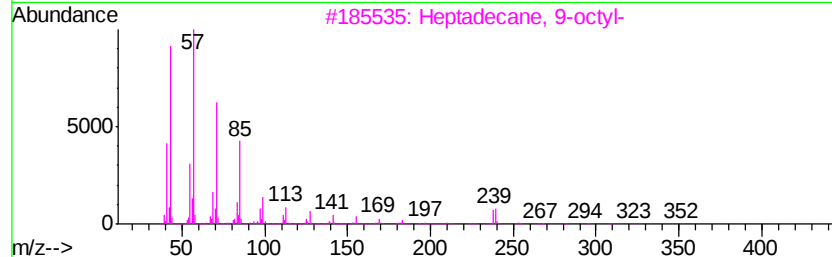
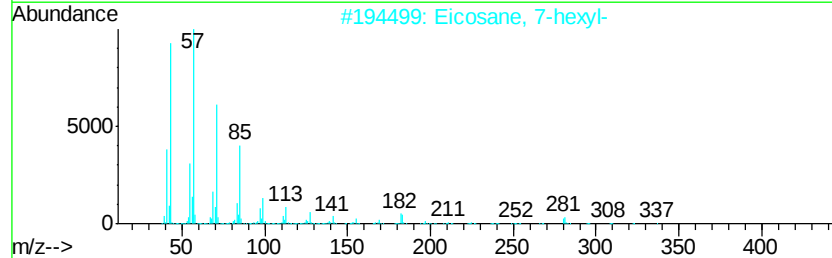
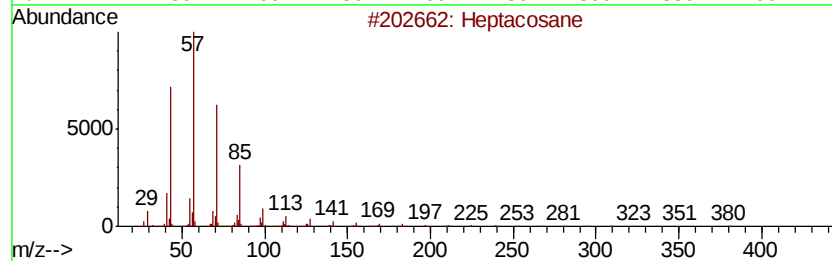
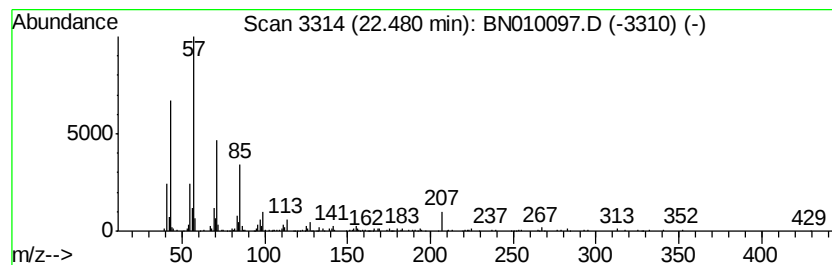
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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.48	4.04 ng/ul	257233	Perylene-d12	23.08

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heptacosane	380	C27H56	000593-49-7	87
2		Eicosane, 7-hexyl-	366	C26H54	055333-99-8	83
3		Heptadecane, 9-octyl-	352	C25H52	007225-64-1	83
4		Eicosane	282	C20H42	000112-95-8	81
5		Hexacosane	366	C26H54	000630-01-3	81



Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN030420\
Data File : BN010097.D
Acq On : 05 Mar 2020 02:27
Operator : CG/JU
Sample : L1617-11
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
DBDS5

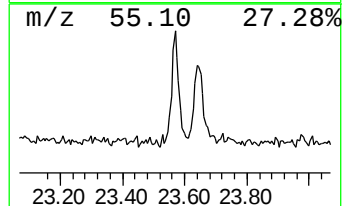
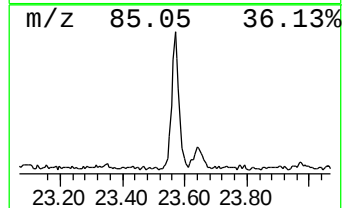
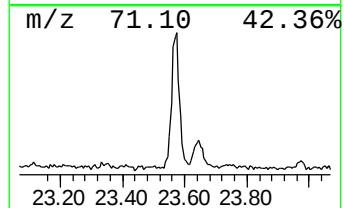
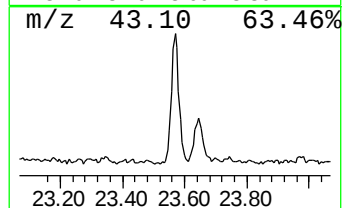
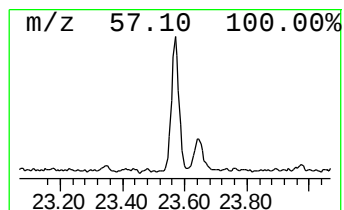
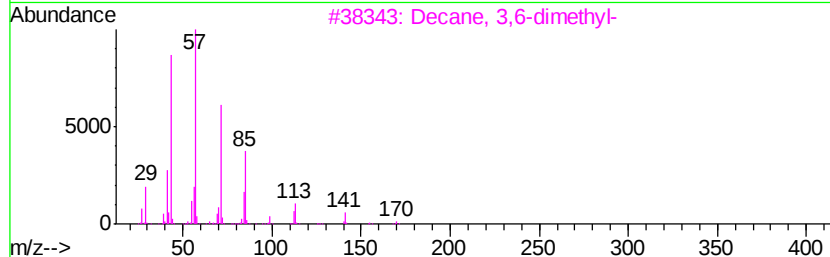
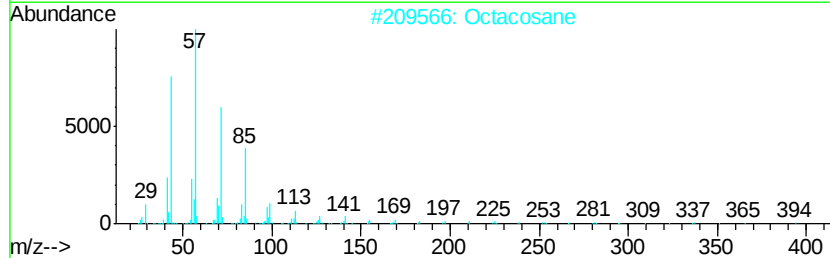
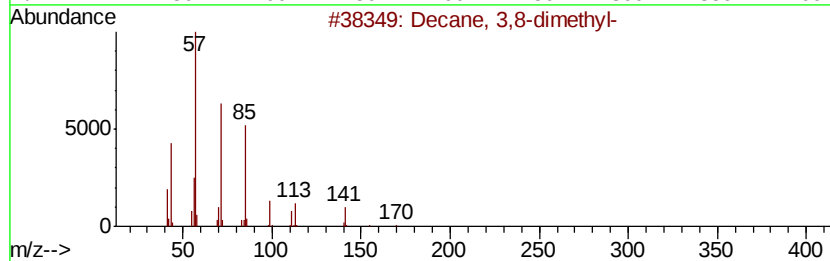
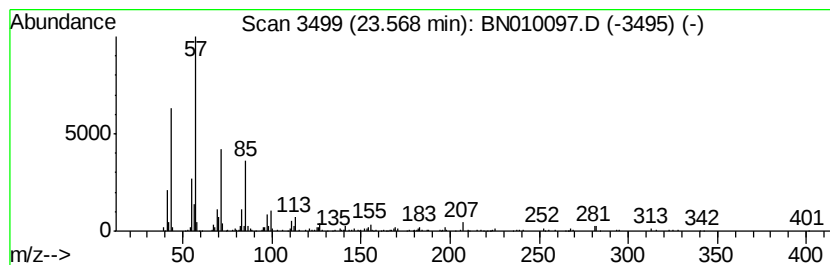
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 (DEL) Alkane: Straight-Chai... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.57	5.51 ng/ul	350904	Perylene-d12	23.08

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Decane, 3,8-dimethyl-	170	C12H26	017312-55-9	83
2			Octacosane	394	C28H58	000630-02-4	81
3			Decane, 3,6-dimethyl-	170	C12H26	017312-53-7	81
4			Hentriacontane	437	C31H64	000630-04-6	76
5			Tridecane, 6-methyl-	198	C14H30	013287-21-3	74



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN030420\
Data File : BN010097.D
Acq On : 05 Mar 2020 02:27
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Sample : L1617-11
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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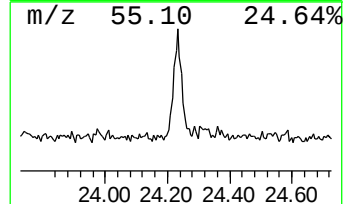
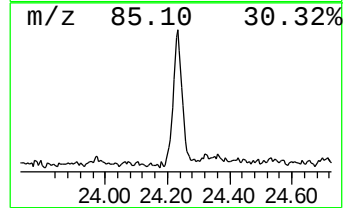
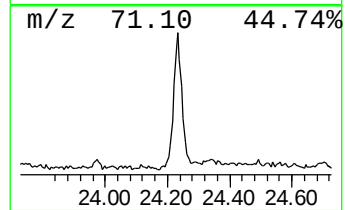
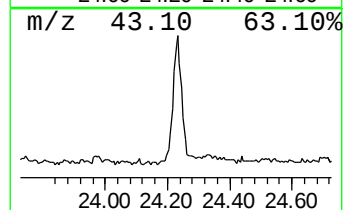
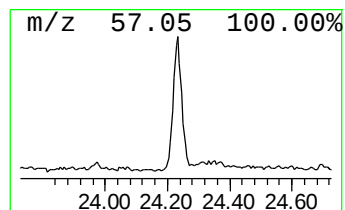
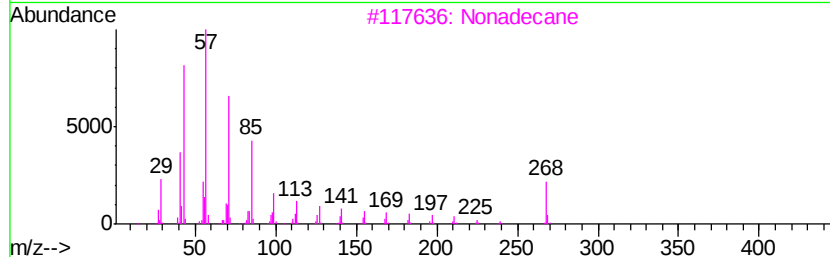
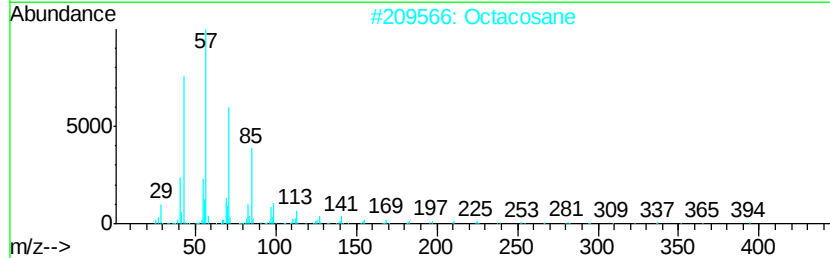
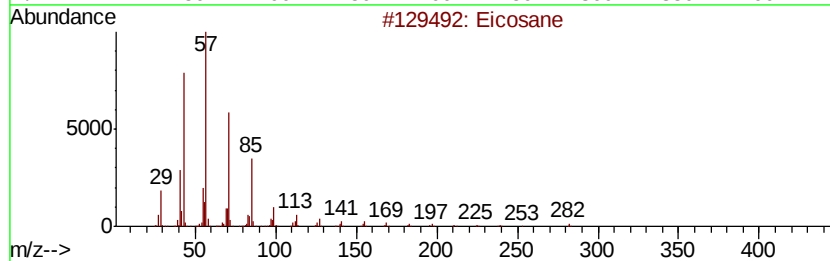
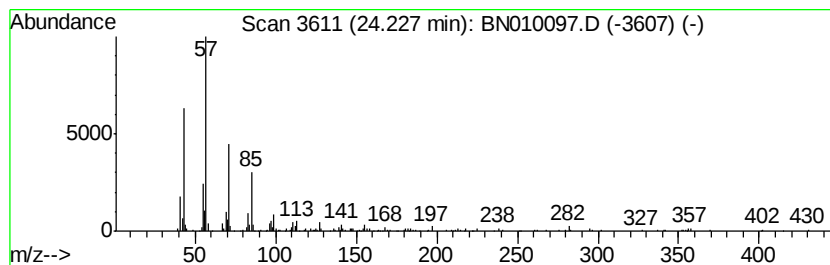
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 (DEL) Alkane: Straight-Chai... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.23	5.58 ng/ul	355251	Perylene-d12	23.08

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Eicosane	282	C20H42	000112-95-8	97
2		Octacosane	394	C28H58	000630-02-4	81
3		Nonadecane	268	C19H40	000629-92-5	81
4		Nonadecane, 9-methyl-	282	C20H42	013287-24-6	74
5		Pentadecane	212	C15H32	000629-62-9	74



Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN030420\
Data File : BN010097.D
Acq On : 05 Mar 2020 02:27
Operator : CG/JU
Sample : L1617-11
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
DBDS5

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
n-Hexadecanoic acid	17.69	4.2	ng/ul	259429	4	16.76	1242150	20.0
(DEL) Alkane: Str...	20.72	11.3	ng/ul	734560	5	20.99	1297380	20.0
1,3-Dioxolane, 2-...	20.79	6.0	ng/ul	386076	5	20.99	1297380	20.0
(DEL) Alkane: Str...	21.59	2.0	ng/ul	130845	5	20.99	1297380	20.0
(DEL) Alkane: Str...	22.02	2.5	ng/ul	162664	5	20.99	1297380	20.0
(DEL) Alkane: Str...	22.48	4.0	ng/ul	257233	6	23.08	1274340	20.0
(DEL) Alkane: Str...	23.57	5.5	ng/ul	350904	6	23.08	1274340	20.0
(DEL) Alkane: Str...	24.23	5.6	ng/ul	355251	6	23.08	1274340	20.0