

Data Path : Z:\svoasrv\HPCHEM1\BNA N\Data\BN030420\
 Data File : BN010085.D
 Acq On : 04 Mar 2020 16:14
 Operator : CG/JU
 Sample : L1641-09
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 CB6Z2

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	3.405	68	71	72	rVV2	2209	2587	0.20%	0.021%
2	3.440	74	77	80	rBV4	2634	2665	0.20%	0.022%
3	3.505	83	88	90	rBV4	2793	4296	0.33%	0.035%
4	3.569	95	99	101	rBV3	2368	3167	0.24%	0.026%
5	3.605	101	105	109	rVB5	4935	6949	0.53%	0.056%
6	3.687	115	119	124	rBV3	8850	11166	0.85%	0.091%
7	3.940	160	162	167	rVB4	2283	3116	0.24%	0.025%
8	4.222	205	210	212	rBV4	2967	4300	0.33%	0.035%
9	4.446	244	248	250	rBV3	2138	2775	0.21%	0.023%
10	4.564	264	268	272	rBV	14242	22032	1.68%	0.179%
11	4.599	272	274	280	rVB4	9497	14050	1.07%	0.114%
12	4.769	300	303	308	rVB5	2117	3008	0.23%	0.024%
13	5.434	410	416	421	rVB3	2995	5415	0.41%	0.044%
14	6.581	605	611	624	rVV2	42298	98910	7.56%	0.804%
15	6.728	631	636	649	rBV	37445	77471	5.92%	0.630%
16	6.910	662	667	679	rBV2	55556	105604	8.07%	0.858%
17	6.999	679	682	686	rVB4	3422	4857	0.37%	0.039%
18	7.357	737	743	757	rVV	276040	492696	37.67%	4.005%
19	8.099	864	869	878	rBV2	38585	76831	5.87%	0.625%
20	8.216	883	889	896	rVV5	3974	8826	0.67%	0.072%
21	8.522	935	941	950	rBV	50322	102861	7.87%	0.836%
22	8.704	966	972	975	rBV5	4373	8478	0.65%	0.069%
23	8.728	975	976	981	rVV2	2991	3926	0.30%	0.032%
24	8.769	981	983	991	rVB3	1885	3326	0.25%	0.027%
25	8.934	1006	1011	1016	rVB	5605	8619	0.66%	0.070%
26	9.234	1056	1062	1075	rVV3	31450	74639	5.71%	0.607%
27	9.781	1150	1155	1163	rBV3	32269	80358	6.14%	0.653%
28	10.104	1201	1210	1226	rBV	371878	680732	52.05%	5.534%
29	10.293	1237	1242	1257	rBV4	30293	93437	7.14%	0.760%
30	11.575	1456	1460	1465	rBV3	2142	2756	0.21%	0.022%
31	12.545	1622	1625	1629	rVB3	2025	2643	0.20%	0.021%
32	12.845	1672	1676	1679	rBV2	2658	3692	0.28%	0.030%
33	13.440	1772	1777	1783	rBV	131893	213235	16.30%	1.733%
34	13.492	1783	1786	1795	rVB	20111	35062	2.68%	0.285%

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35	13.692	1813	1820	1830	rBV	141335	223134	17.06%	1.814%
36	13.922	1857	1859	1860	rVV2	3356	2605	0.20%	0.021%
37	13.945	1860	1863	1867	rVV3	4162	6879	0.53%	0.056%
38	14.004	1867	1873	1882	rVV	617530	939862	71.86%	7.640%
39	14.069	1882	1884	1887	rVV3	4903	5805	0.44%	0.047%
40	14.316	1921	1926	1932	rBV7	11741	29385	2.25%	0.239%
41	14.839	2010	2015	2019	rBV2	2833	4106	0.31%	0.033%
42	14.910	2024	2027	2030	rBV3	3068	3840	0.29%	0.031%
43	15.016	2040	2045	2053	rBV	187239	302593	23.14%	2.460%
44	15.204	2071	2077	2085	rBV7	19159	45311	3.46%	0.368%
45	15.257	2085	2086	2087	rVV	5033	2512	0.19%	0.020%
46	15.275	2087	2089	2092	rVV2	3567	3758	0.29%	0.031%
47	15.298	2092	2093	2098	rVB2	2967	3046	0.23%	0.025%
48	15.610	2141	2146	2147	rBV2	2066	2696	0.21%	0.022%
49	15.939	2197	2202	2204	rBV3	2306	3918	0.30%	0.032%
50	16.339	2266	2270	2274	rVV4	4065	6679	0.51%	0.054%
51	16.563	2303	2308	2313	rBV4	4833	8388	0.64%	0.068%
52	16.598	2313	2314	2318	rVV4	3816	3206	0.25%	0.026%
53	16.763	2335	2342	2348	rBV	772345	1133218	86.65%	9.212%
54	16.869	2354	2360	2374	rVB2	201798	330683	25.28%	2.688%
55	17.181	2409	2413	2416	rBV4	3488	4865	0.37%	0.040%
56	17.633	2486	2490	2491	rBV3	2525	3028	0.23%	0.025%
57	17.657	2492	2494	2498	rVB2	3251	3859	0.30%	0.031%
58	17.710	2498	2503	2509	rBV6	7672	14130	1.08%	0.115%
59	17.845	2522	2526	2531	rBV5	6427	10783	0.82%	0.088%
60	17.892	2531	2534	2538	rVB5	3494	4878	0.37%	0.040%
61	17.986	2549	2550	2555	rVB3	2710	3286	0.25%	0.027%
62	18.128	2571	2574	2577	rBV4	5468	6757	0.52%	0.055%
63	18.169	2579	2581	2585	rVB3	5170	4174	0.32%	0.034%
64	18.398	2617	2620	2622	rBV3	3741	3785	0.29%	0.031%
65	18.422	2622	2624	2627	rVB3	2568	2543	0.19%	0.021%
66	18.463	2627	2631	2633	rBV4	2581	3577	0.27%	0.029%
67	18.563	2646	2648	2650	rBV2	3278	2965	0.23%	0.024%
68	18.639	2657	2661	2665	rBV5	7553	9729	0.74%	0.079%
69	18.722	2674	2675	2681	rVB4	4174	5355	0.41%	0.044%
70	18.845	2691	2696	2702	rVV	58160	81643	6.24%	0.664%
71	19.004	2719	2723	2725	rBV4	4079	5638	0.43%	0.046%

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72	19.139	2742	2746	2748	rBV3	5841	9589	0.73%	0.078%
73	19.180	2748	2753	2757	rBV	275182	418149	31.97%	3.399%
74	19.263	2765	2767	2768	rBV2	3118	2625	0.20%	0.021%
75	19.304	2770	2774	2776	rBV5	4669	6054	0.46%	0.049%
76	19.357	2782	2783	2785	rBV2	4544	3631	0.28%	0.030%
77	19.692	2838	2840	2842	rBV3	5296	4334	0.33%	0.035%
78	19.733	2843	2847	2850	rVV5	14839	24032	1.84%	0.195%
79	19.769	2850	2853	2860	rVB3	21412	27833	2.13%	0.226%
80	19.880	2870	2872	2876	rVB4	7455	7770	0.59%	0.063%
81	19.945	2881	2883	2886	rVV4	5701	5371	0.41%	0.044%
82	20.033	2894	2898	2901	rBV6	5176	9300	0.71%	0.076%
83	20.269	2935	2938	2941	rBV4	16325	18190	1.39%	0.148%
84	20.686	3003	3009	3012	rBV	116302	189270	14.47%	1.539%
85	20.733	3012	3017	3022	rVV2	251007	389621	29.79%	3.167%
86	20.786	3022	3026	3035	rVV	479259	699284	53.47%	5.684%
87	20.857	3036	3038	3041	rVB2	24590	29237	2.24%	0.238%
88	20.992	3055	3061	3066	rBV	992175	1307824	100.00%	10.631%
89	21.116	3079	3082	3085	rBV2	43997	54192	4.14%	0.441%
90	21.174	3089	3092	3096	rVB	53043	57518	4.40%	0.468%
91	21.598	3160	3164	3173	rBV2	173166	304989	23.32%	2.479%
92	22.021	3233	3236	3244	rVB	98807	132527	10.13%	1.077%
93	22.486	3311	3315	3319	rBV2	289629	401853	30.73%	3.267%
94	22.521	3319	3321	3328	rVB2	108996	153327	11.72%	1.246%
95	22.957	3391	3395	3399	rBV	209655	322647	24.67%	2.623%
96	22.998	3399	3402	3409	rVB	165987	252928	19.34%	2.056%
97	23.086	3411	3417	3429	rVB	684012	1276919	97.64%	10.380%
98	23.280	3447	3450	3455	rVB4	103013	138074	10.56%	1.122%
99	23.580	3496	3501	3508	rVB	201066	356429	27.25%	2.897%
100	24.239	3609	3613	3620	rVB	103455	201216	15.39%	1.636%

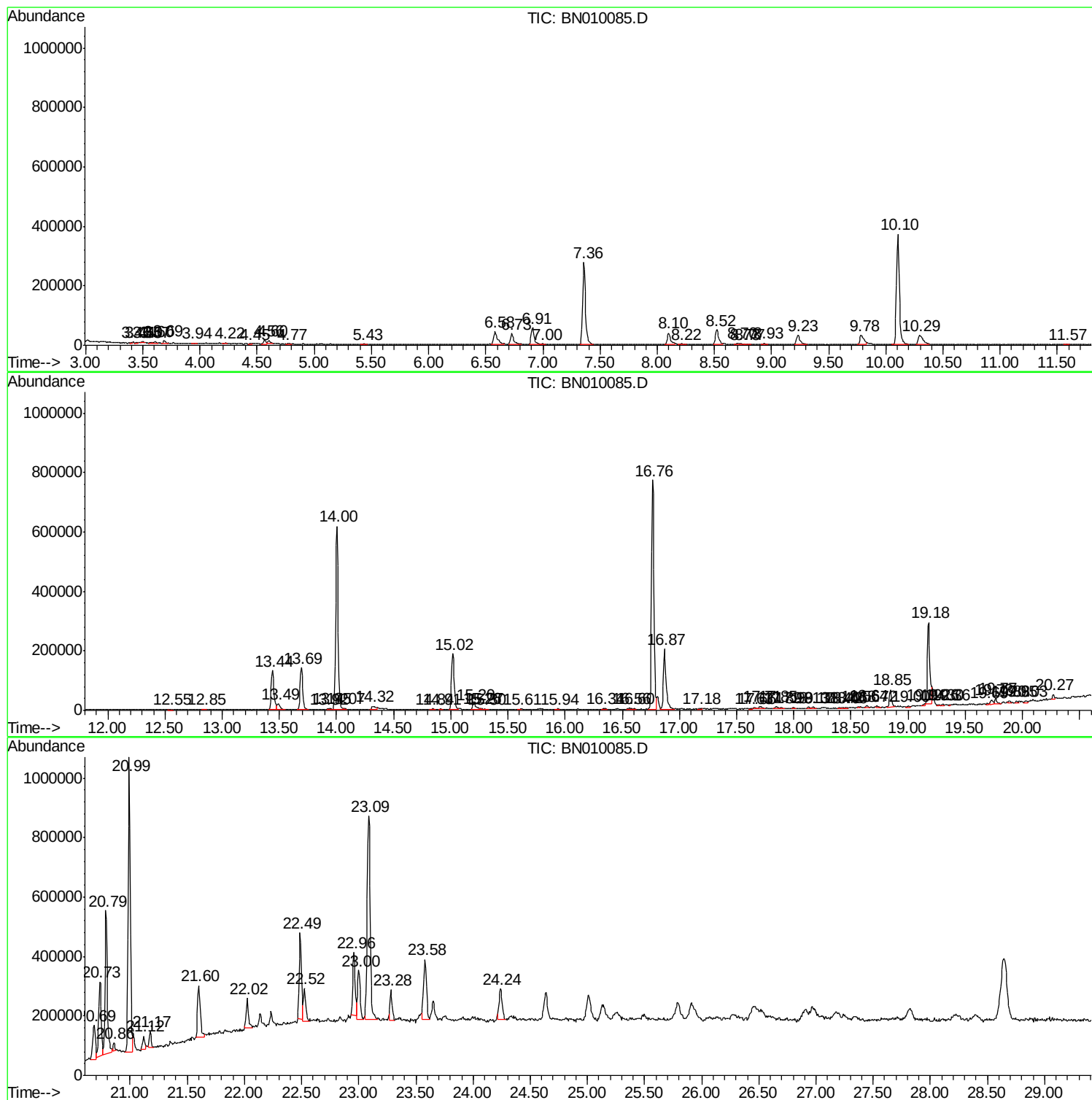
Sum of corrected areas: 12301837

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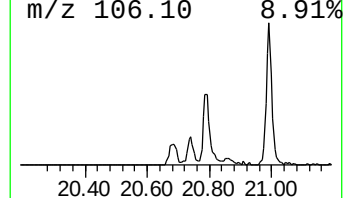
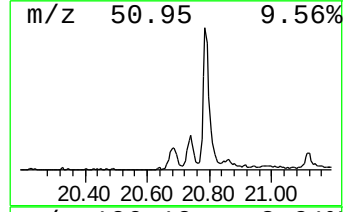
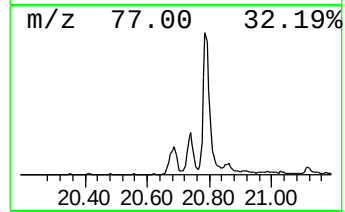
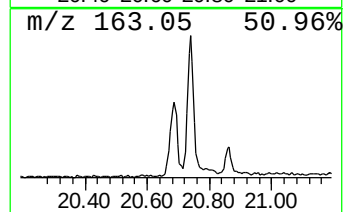
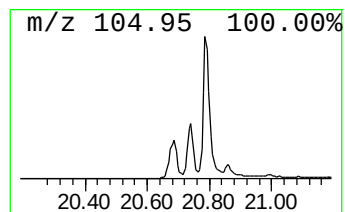
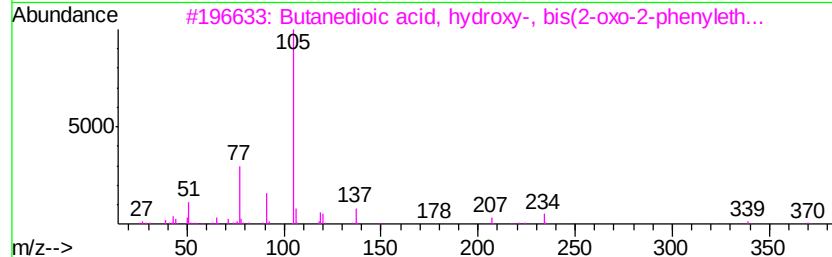
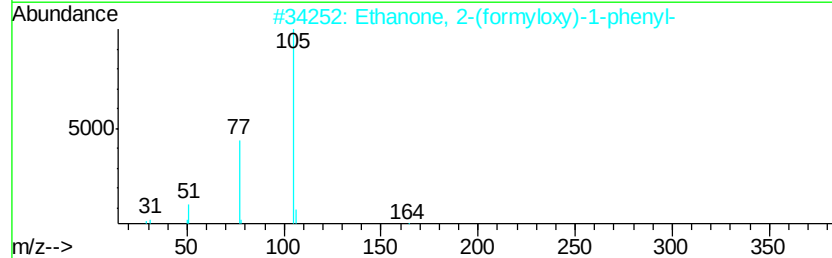
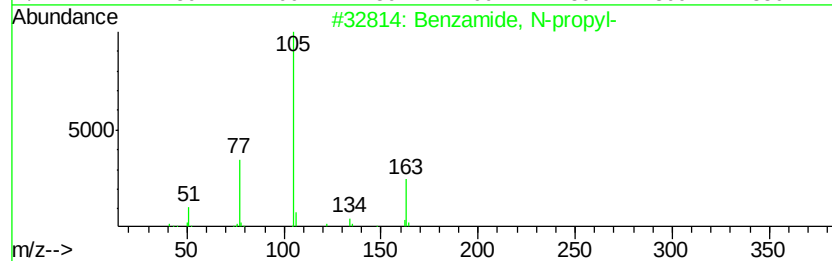
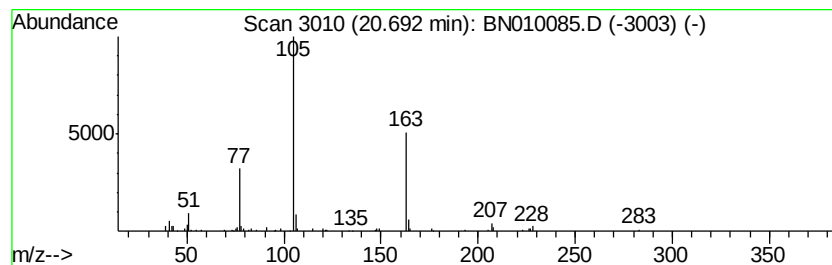
Instrument :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN021220MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Benzamide, N-propyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.		
20.69	2.89 ng/ul	189270	Chrysene-d12	20.99		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Benzamide, N-propyl-	163	C10H13NO	010546-70-0	80
2		Ethanone, 2-(formyloxy)-1-phenyl-	164	C9H8O3	055153-12-3	38
3		Butanedioic acid, hydroxy-, bis(...	370	C20H18O7	058265-78-4	38
4		2-Methoxyphenyl benzoate	228	C14H12O3	000531-37-3	38
5		Pentanediamide, N,N'-di-benzoyloxy-	370	C19H18N2O6	1000253-26-3	35



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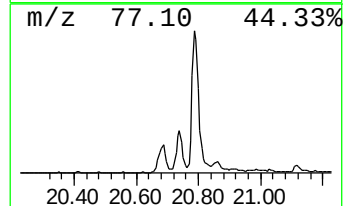
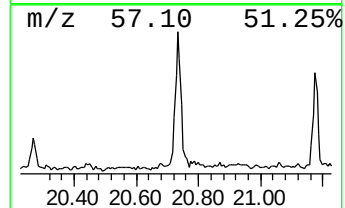
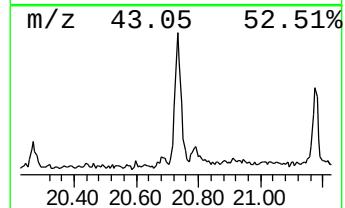
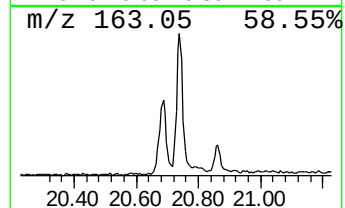
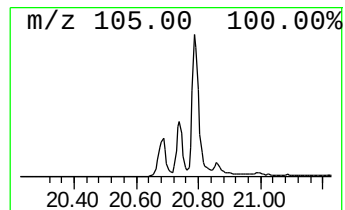
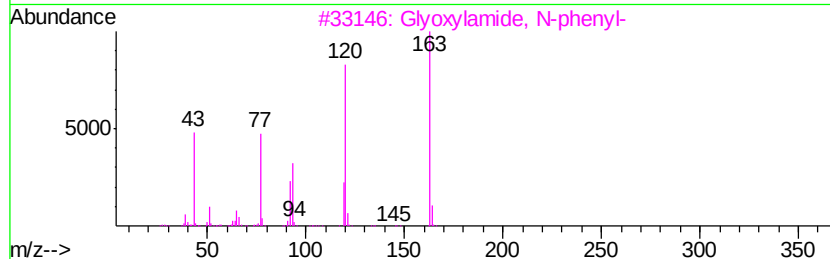
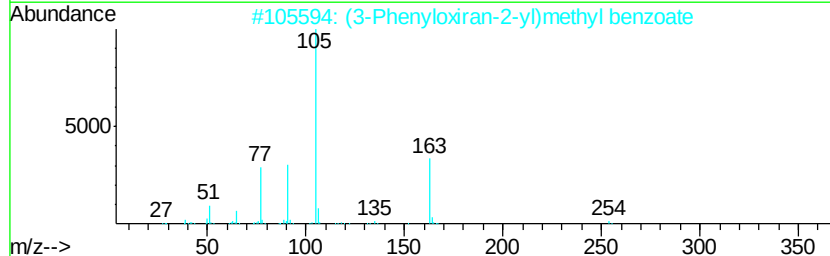
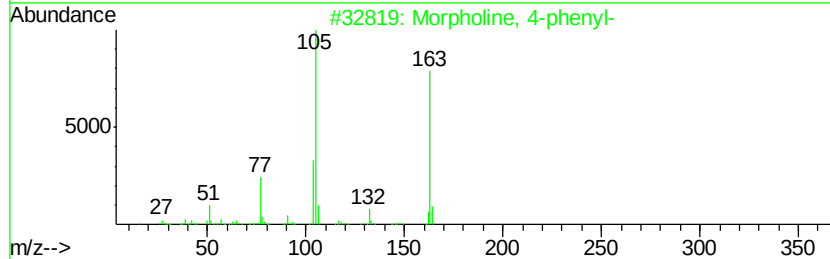
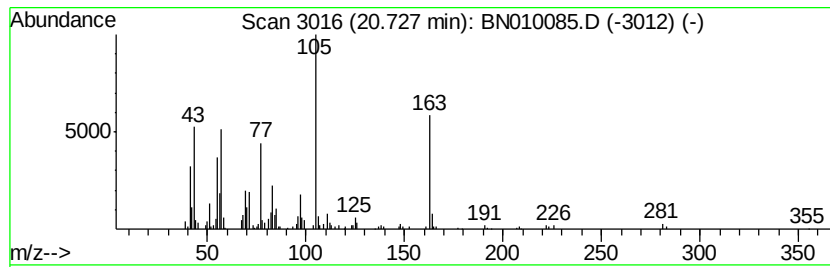
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Quant Title : SVOA CALIBRATION

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

Peak Number 2 Morpholine, 4-phenyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD		R.T.	
20.73	5.96 ng/ul	389621	Chrysene-d12		20.99	
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Morpholine, 4-phenyl-	163	C10H13NO	000092-53-5	52
2		(3-Phenyloxiran-2-yl)methyl benz...	254	C16H14O3	1000366-96-1	47
3		Glvoxylamide, N-phenyl-	163	C9H9NO2	032331-52-5	43
4		Benzoic acid, 4-methylpent-2-yl ...	206	C13H18O2	1000367-94-2	35
5		Benzo[b]thiophene-3-carboxylic a...	343	C18H17NO4S	096334-43-9	22



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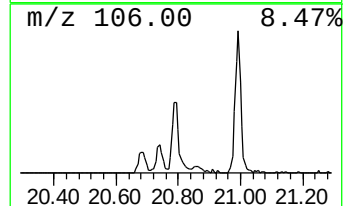
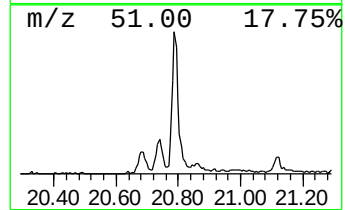
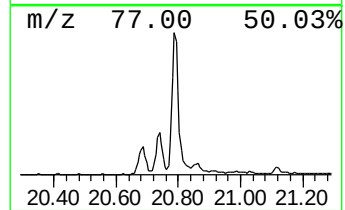
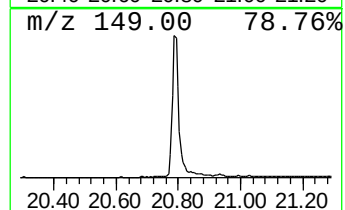
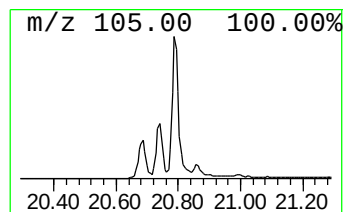
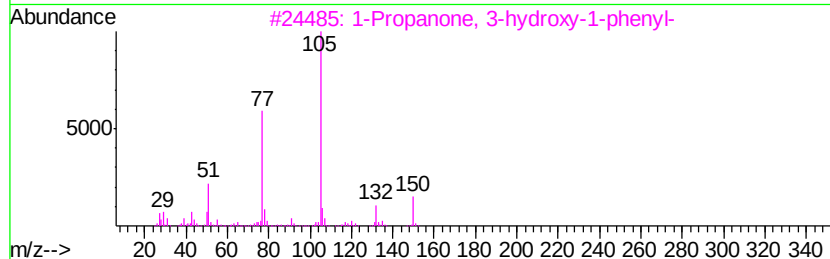
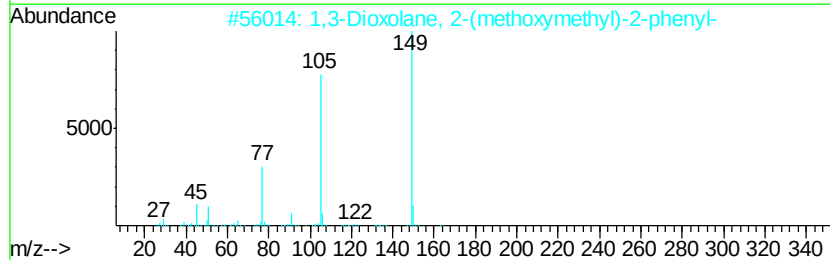
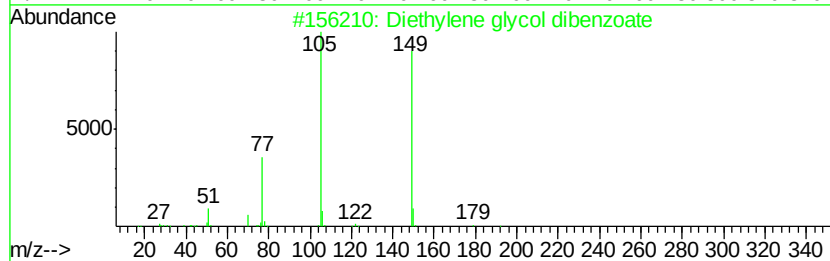
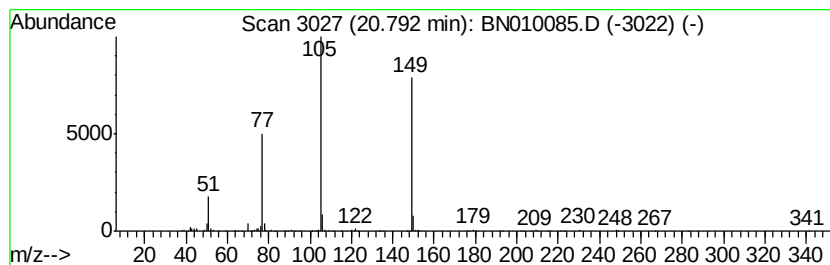
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 Diethylene glycol dibenzoate Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.		
20.79	10.69 ng/ul	699284	Chrysene-d12	20.99		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Diethylene glycol dibenzoate	314	C18H18O5	000120-55-8	78
2		1,3-Dioxolane, 2-(methoxymethyl)...	194	C11H14O3	1000156-71-2	53
3		1-Propanone, 3-hydroxy-1-phenyl-	150	C9H10O2	005650-41-9	47
4		4'-Nitrobenzanilide	242	C13H10N2O3	003393-96-2	43
5		1,3,4-Oxadiazin-6-one, 5-isoprop...	216	C12H12N2O2	173973-78-9	43



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Data File : BN010085.D
Acq On : 04 Mar 2020 16:14
Operator : CG/JU
Sample : L1641-09
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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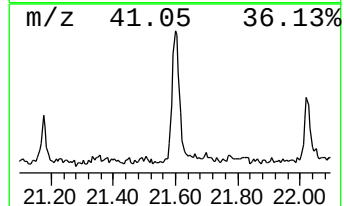
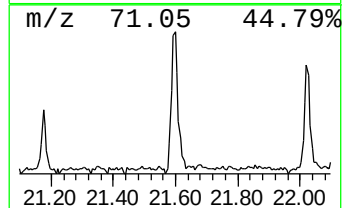
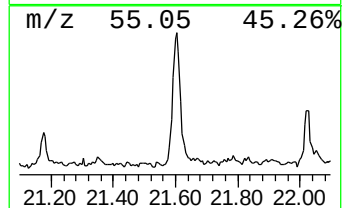
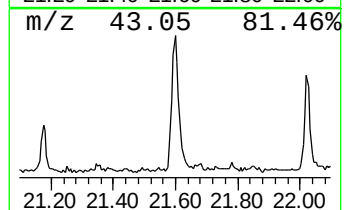
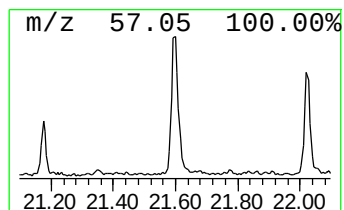
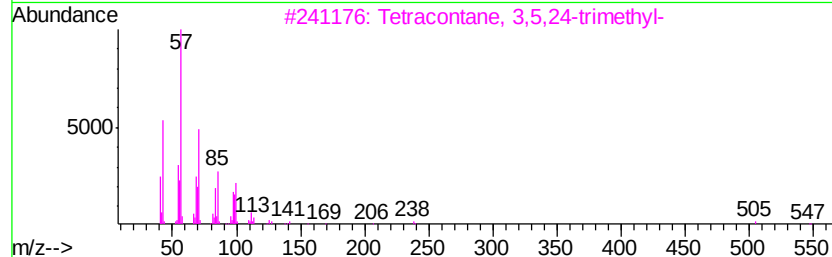
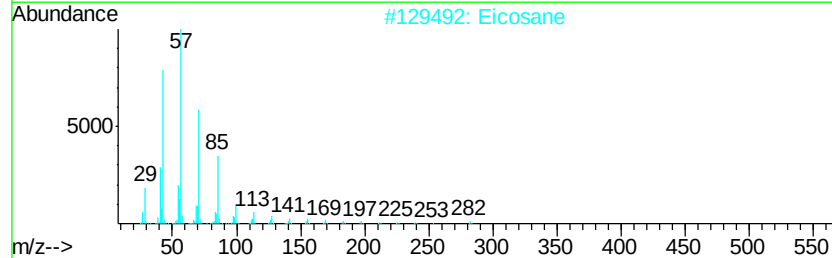
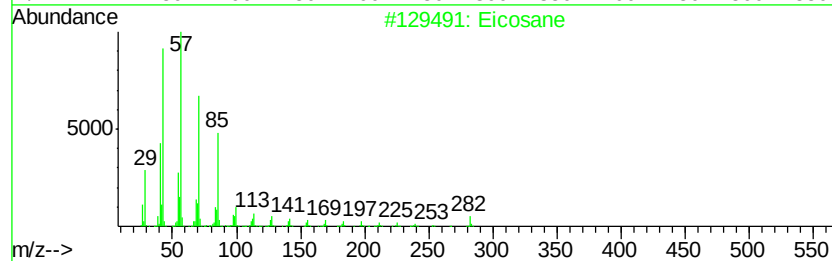
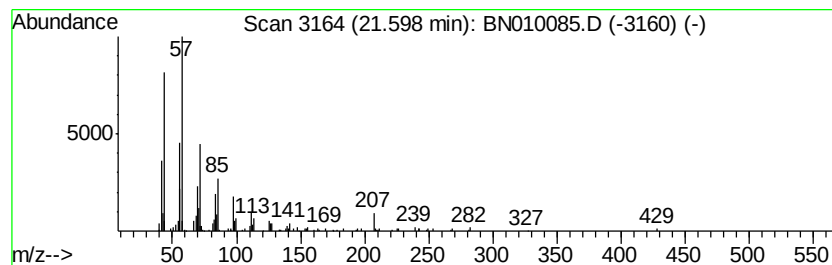
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 (DEL) Alkane: Straight-Chai... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.60	4.66 ng/ul	304989	Chrysene-d12	20.99

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Eicosane	282	C20H42	000112-95-8	91
2		Eicosane	282	C20H42	000112-95-8	90
3		Tetracontane, 3,5,24-trimethyl-	605	C43H88	055162-61-3	86
4		Sulfurous acid, butyl octadecyl ...	390	C22H46O3S	1000309-18-5	86
5		Hexadecane, 1-chloro-	260	C16H33Cl	004860-03-1	80



Data Path : Z:\svoasrv\HPCHEM1\BNA N\Data\BN030420\
Data File : BN010085.D
Acq On : 04 Mar 2020 16:14
Operator : CG/JU
Sample : L1641-09
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
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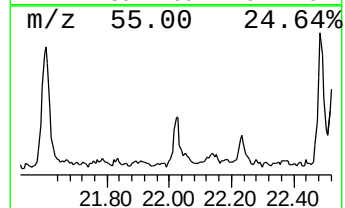
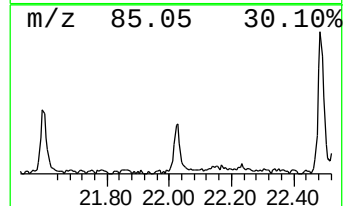
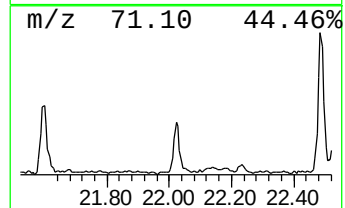
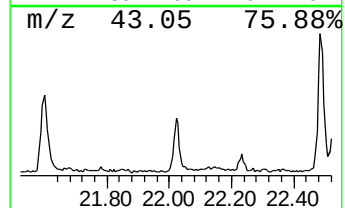
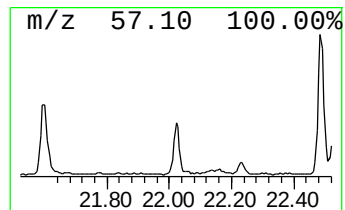
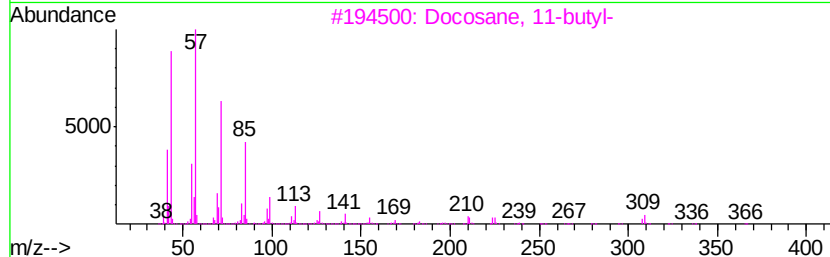
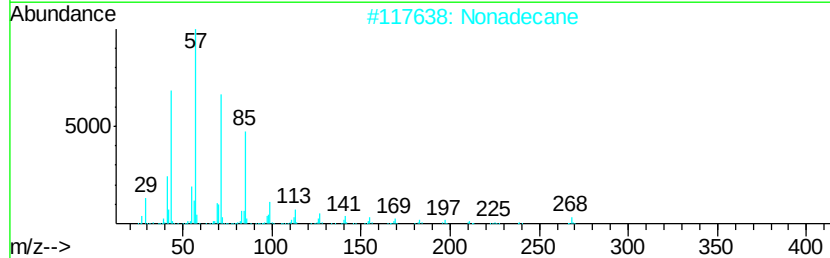
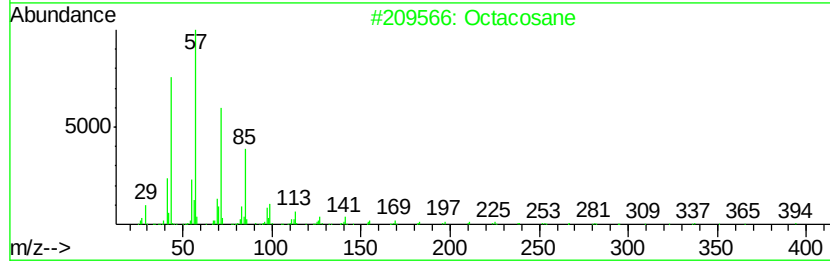
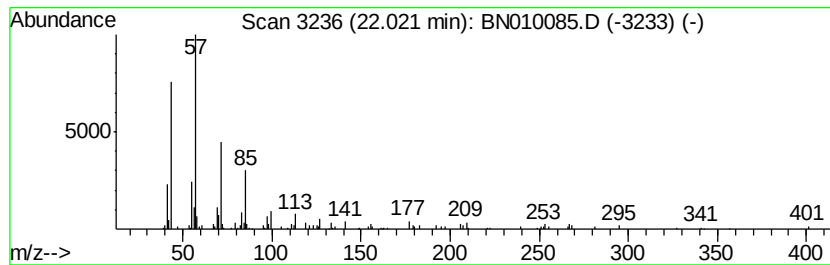
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 (DEL) Alkane: Straight-Chai... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.02	2.03 ng/ul	132527	Chrysene-d12	20.99

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octacosane	394	C28H58	000630-02-4	87
2		Nonadecane	268	C19H40	000629-92-5	76
3		Docosane, 11-butyl-	366	C26H54	013475-76-8	68
4		Hexacosane	366	C26H54	000630-01-3	68
5		Heptadecane, 9-octyl-	352	C25H52	007225-64-1	68



Data Path : Z:\svoasrv\HPCHEM1\BNA N\Data\BN030420\
Data File : BN010085.D
Acq On : 04 Mar 2020 16:14
Operator : CG/JU
Sample : L1641-09
Misc :
ALS Vial : 8 Sample Multiplier: 1

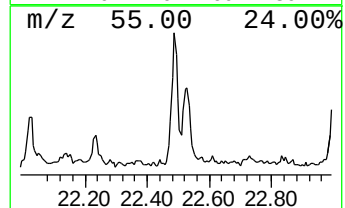
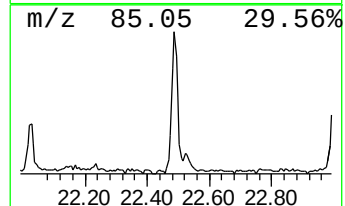
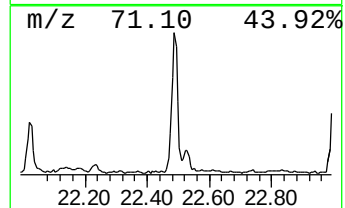
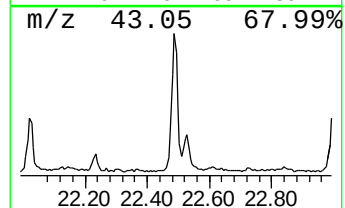
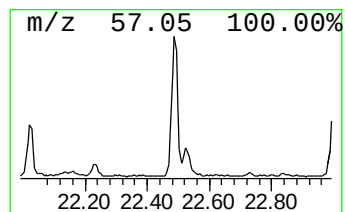
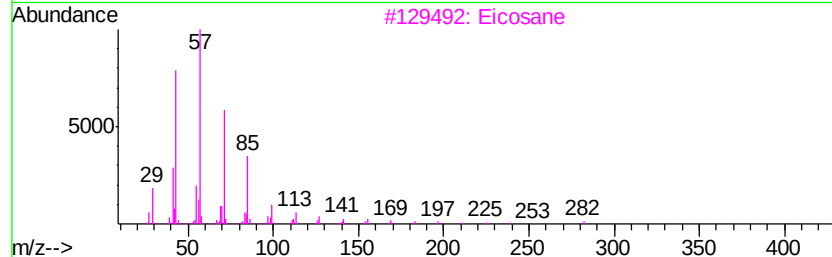
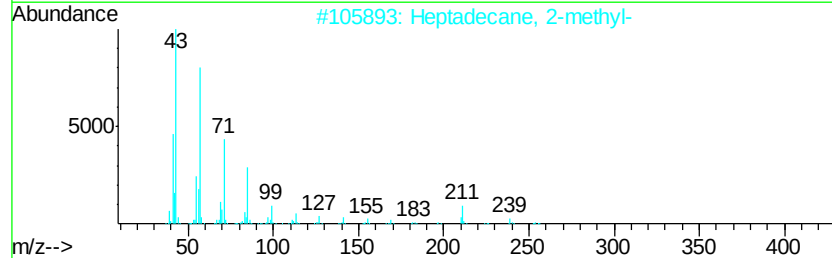
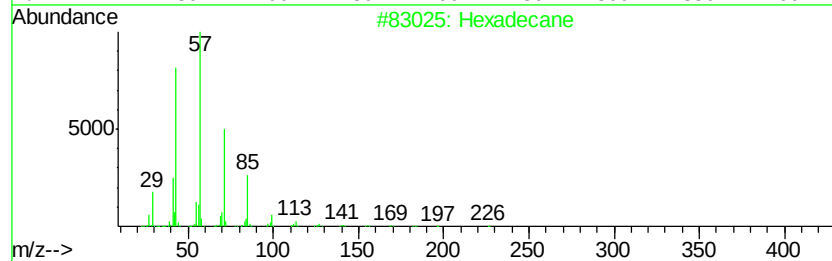
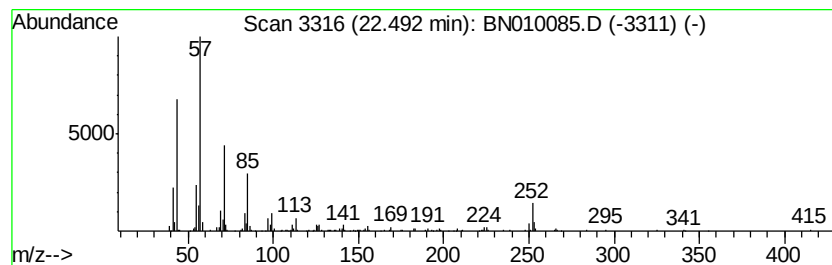
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ClientSampleId :
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 (DEL) Alkane: Straight-Chai... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.49	6.29 ng/ul	401853	Perylene-d12	23.09
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Hexadecane	226 C16H34	000544-76-3	93
2	Heptadecane, 2-methyl-	254 C18H38	001560-89-0	70
3	Eicosane	282 C20H42	000112-95-8	68
4	Hexatriacontane	507 C36H74	000630-06-8	68
5	Sulfurous acid, butyl undecyl ester	292 C15H32O3S	1000309-17-8	64



Data Path : Z:\svoasrv\HPCHEM1\BNA N\Data\BN030420\
Data File : BN010085.D
Acq On : 04 Mar 2020 16:14
Operator : CG/JU
Sample : L1641-09
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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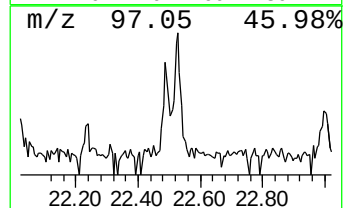
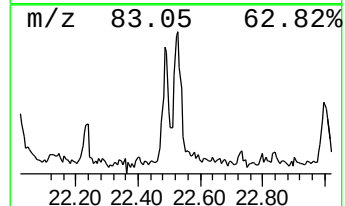
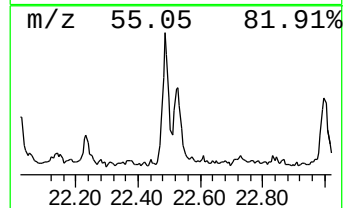
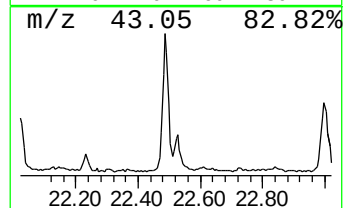
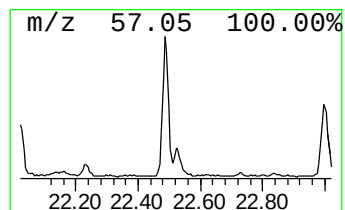
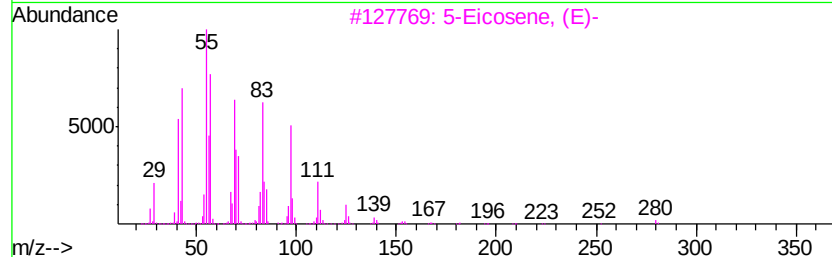
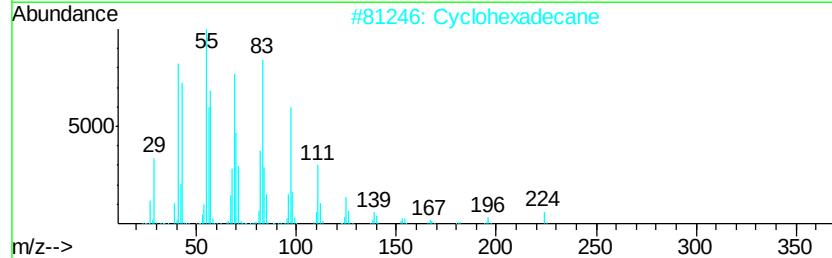
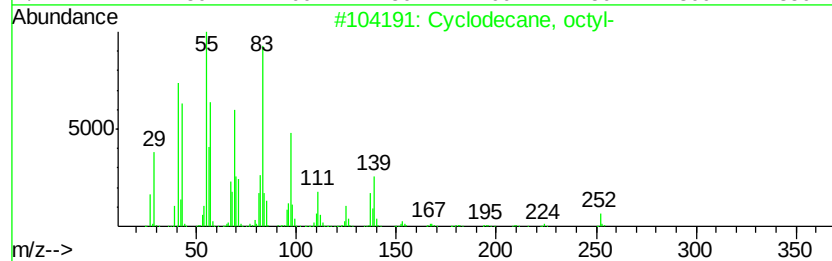
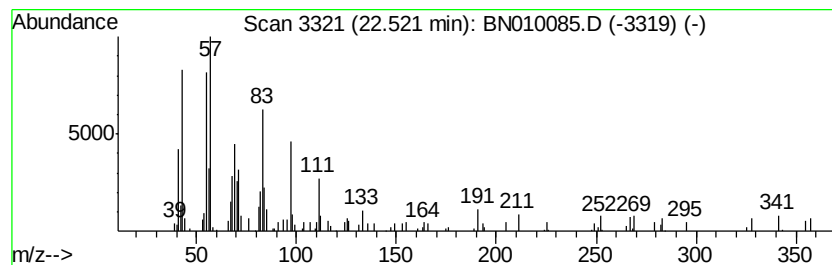
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 (DEL) Alkane: Cyclic22.52 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.52	2.40 ng/ul	153327	Perylene-d12	23.09

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclodecane, octyl-	252	C18H36	016539-09-6	83
2		Cyclohexadecane	224	C16H32	000295-65-8	58
3		5-Eicosene, (E)-	280	C20H40	074685-30-6	58
4		3-Chloropropionic acid, heptadec...	346	C20H39ClO2	1000283-05-1	58
5		cis-13-Octadecenoic acid	282	C18H34O2	013126-39-1	58



Data Path : Z:\svoasrv\HPCHEM1\BNA N\Data\BN030420\
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Acq On : 04 Mar 2020 16:14
Operator : CG/JU
Sample : L1641-09
Misc :
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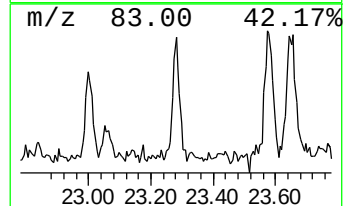
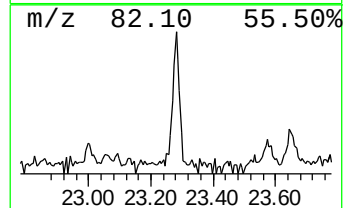
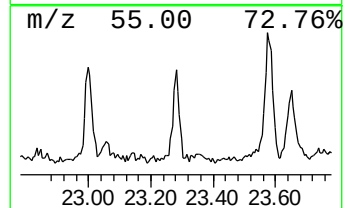
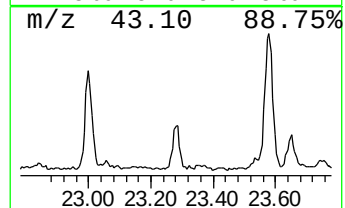
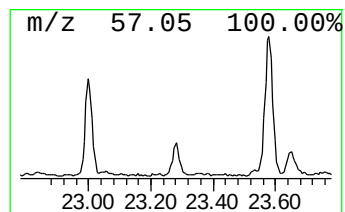
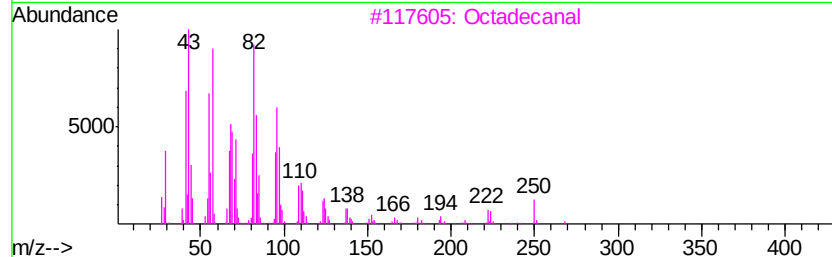
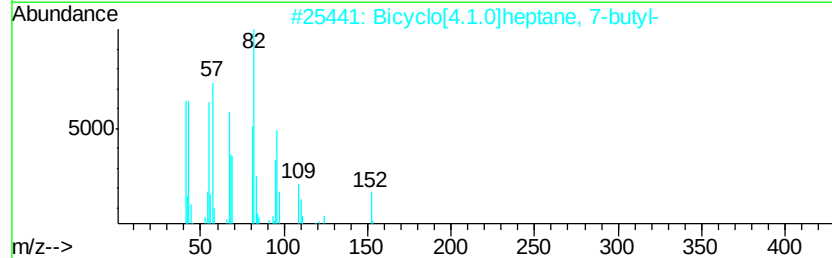
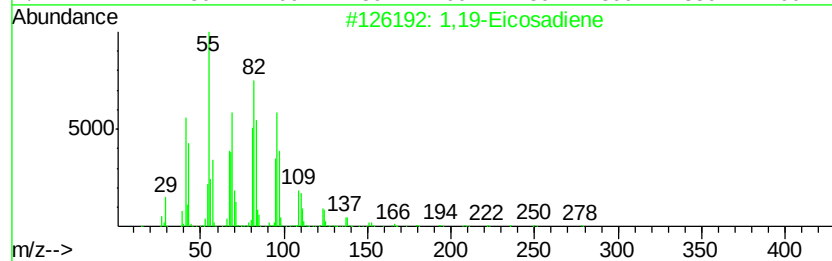
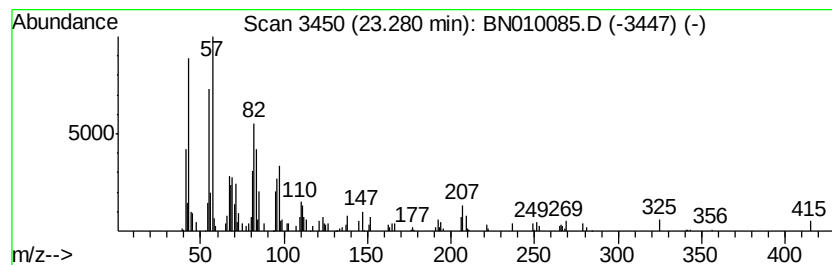
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 1,19-Eicosadiene Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.28	2.16 ng/ul	138074	Perylene-d12	23.09
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	1,19-Eicosadiene		278 C20H38	014811-95-1 58
2	Bicyclo[4.1.0]heptane, 7-butyl-		152 C11H20	018645-10-8 55
3	Octadecanal		268 C18H36O	000638-66-4 52
4	Hexadecane, 1-(ethenyl)-		268 C18H36O	000822-28-6 43
5	Z,Z-3,13-Octadecadien-1-ol		266 C18H34O	1000131-10-7 41



Data Path : Z:\svoasrv\HPCHEM1\BNA N\Data\BN030420\
Data File : BN010085.D
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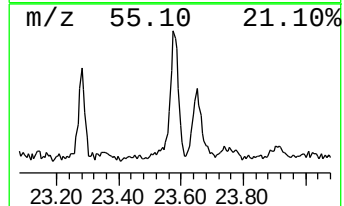
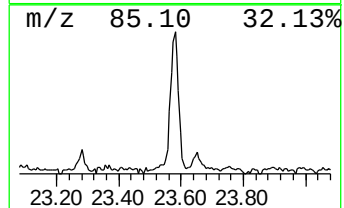
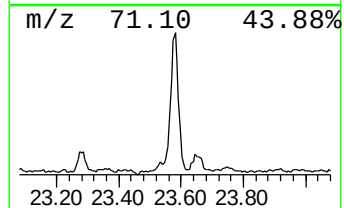
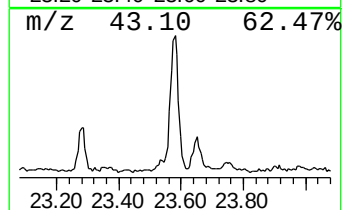
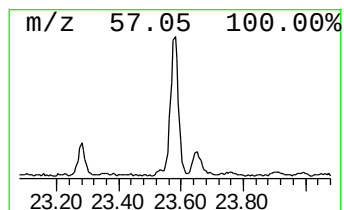
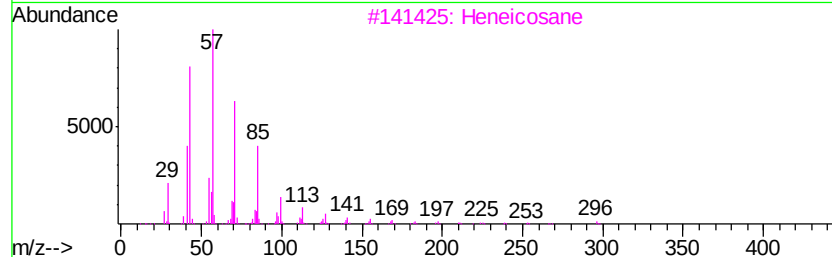
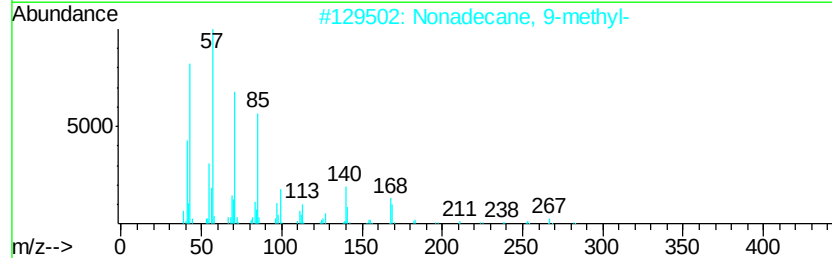
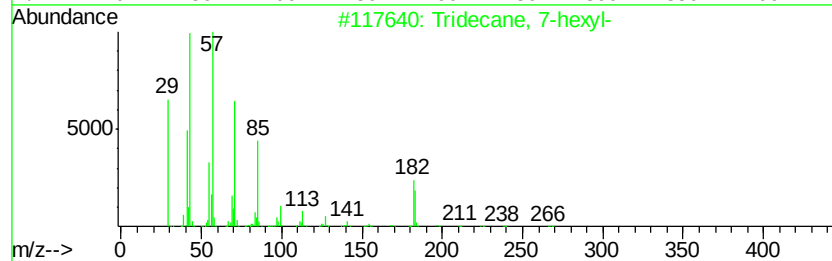
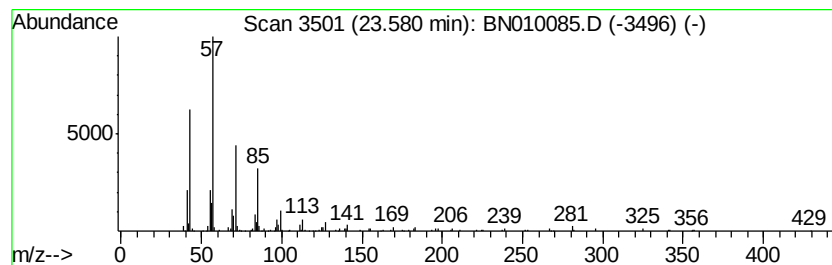
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 (DEL) Alkane: Straight-Chai... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.58	5.58 ng/ul	356429	Perylene-d12	23.09

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tridecane, 7-hexyl-	268	C19H40	007225-66-3	91
2		Nonadecane, 9-methyl-	282	C20H42	013287-24-6	89
3		Heneicosane	296	C21H44	000629-94-7	87
4		Eicosane, 7-hexyl-	366	C26H54	055333-99-8	83
5		Hexatriacontane	507	C36H74	000630-06-8	83



Data Path : Z:\svoasrv\HPCHEM1\BNA N\Data\BN030420\
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Misc :
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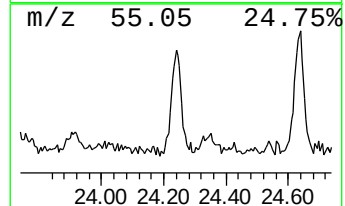
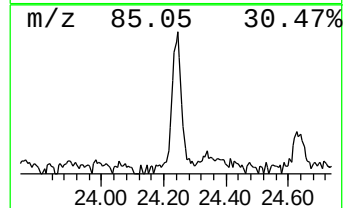
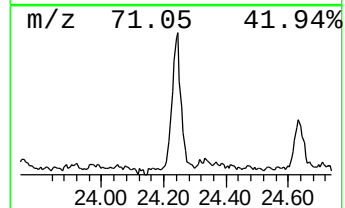
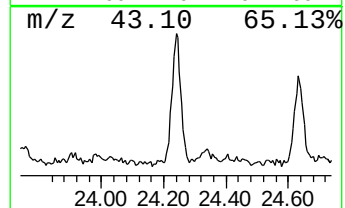
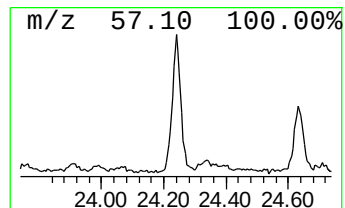
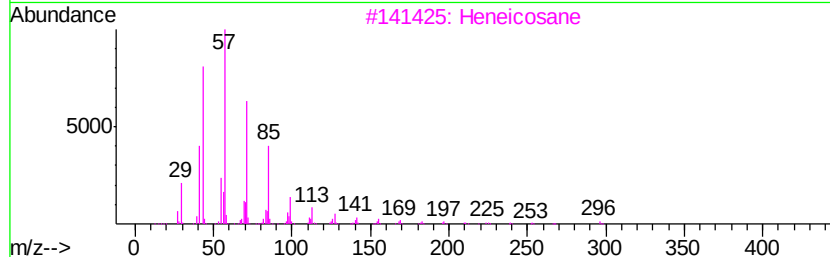
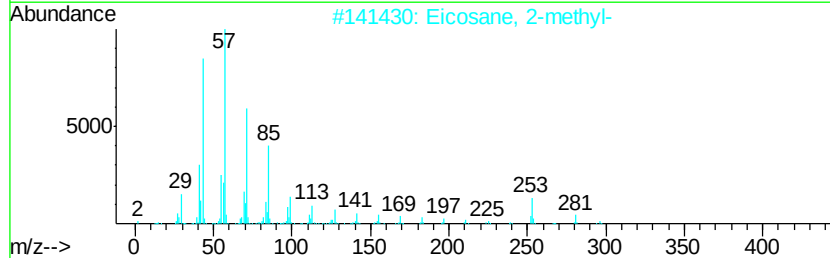
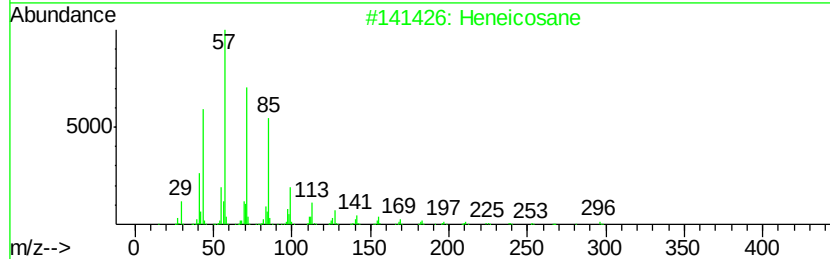
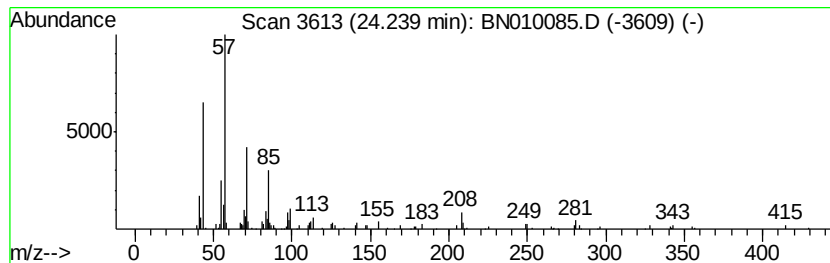
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 (DEL) Alkane: Straight-Chai... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.24	3.15 ng/ul	201216	Perylene-d12	23.09

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heneicosane	296	C21H44	000629-94-7	93
2		Eicosane, 2-methyl-	296	C21H44	001560-84-5	89
3		Heneicosane	296	C21H44	000629-94-7	86
4		Octacosane	394	C28H58	000630-02-4	81
5		Eicosane, 7-hexyl-	366	C26H54	055333-99-8	76



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN030420\
 Data File : BN010085.D
 Acq On : 04 Mar 2020 16:14
 Operator : CG/JU
 Sample : L1641-09
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 CB6Z2

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
Benzamide, N-propyl-	20.69	2.9	ng/ul	189270	5	20.99	1307820	20.0
Morpholine, 4-phe...	20.73	6.0	ng/ul	389621	5	20.99	1307820	20.0
Diethylene glycol...	20.79	10.7	ng/ul	699284	5	20.99	1307820	20.0
(DEL) Alkane: Str...	21.60	4.7	ng/ul	304989	5	20.99	1307820	20.0
(DEL) Alkane: Str...	22.02	2.0	ng/ul	132527	5	20.99	1307820	20.0
(DEL) Alkane: Str...	22.49	6.3	ng/ul	401853	6	23.09	1276920	20.0
(DEL) Alkane: Cyc...	22.52	2.4	ng/ul	153327	6	23.09	1276920	20.0
1,19-Eicosadiene	23.28	2.2	ng/ul	138074	6	23.09	1276920	20.0
(DEL) Alkane: Str...	23.58	5.6	ng/ul	356429	6	23.09	1276920	20.0
(DEL) Alkane: Str...	24.24	3.1	ng/ul	201216	6	23.09	1276920	20.0