

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN022220\
 Data File : BN009817.D
 Acq On : 21 Feb 2020 20:55
 Operator : CG/JU
 Sample : L1535-03
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 CB6Q1

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	3.040	7	9	19	rVB7	9315	19649	1.06%	0.108%
2	3.640	105	111	115	rVB3	12459	18695	1.01%	0.103%
3	3.722	121	125	130	rBV4	8120	13974	0.75%	0.077%
4	3.940	157	162	166	rBV	38991	54691	2.95%	0.301%
5	4.199	202	206	210	rBV5	4476	7660	0.41%	0.042%
6	4.252	211	215	217	rVB3	4775	5693	0.31%	0.031%
7	4.328	224	228	236	rBV	30106	47141	2.54%	0.260%
8	4.599	269	274	282	rBV2	26126	42051	2.27%	0.232%
9	5.463	419	421	426	rVB4	2765	4432	0.24%	0.024%
10	6.605	609	615	629	rVV2	72181	153940	8.29%	0.848%
11	6.757	634	641	650	rBV2	64385	113431	6.11%	0.625%
12	6.940	666	672	681	rBV2	90655	153169	8.25%	0.844%
13	7.034	685	688	692	rBV4	4560	6355	0.34%	0.035%
14	7.393	743	749	758	rBV	450262	727972	39.22%	4.011%
15	7.493	763	766	771	rVB4	2951	4046	0.22%	0.022%
16	8.116	867	872	882	rBV2	80299	154783	8.34%	0.853%
17	8.234	888	892	894	rBV3	3860	4696	0.25%	0.026%
18	8.546	938	945	953	rBV	81540	154995	8.35%	0.854%
19	8.728	972	976	980	rBV3	8424	11350	0.61%	0.063%
20	8.875	994	1001	1005	rBV3	4173	8501	0.46%	0.047%
21	8.975	1014	1018	1025	rVB3	5655	9119	0.49%	0.050%
22	9.251	1059	1065	1080	rBV	59500	123147	6.64%	0.679%
23	9.798	1152	1158	1176	rBV2	71868	164680	8.87%	0.907%
24	10.140	1208	1216	1231	rBV	642278	1080184	58.20%	5.952%
25	10.310	1239	1245	1256	rBV	63287	161959	8.73%	0.892%
26	12.881	1678	1682	1686	rVB3	3255	4124	0.22%	0.023%
27	13.351	1757	1762	1766	rVB3	3412	4708	0.25%	0.026%
28	13.463	1774	1781	1787	rBV	193145	293015	15.79%	1.615%
29	13.510	1787	1789	1797	rVB	16828	24345	1.31%	0.134%
30	13.722	1818	1825	1832	rBV	222670	327602	17.65%	1.805%
31	13.945	1858	1863	1864	rBV3	3601	4399	0.24%	0.024%
32	14.034	1872	1878	1885	rBV	937410	1392735	75.04%	7.674%
33	14.098	1885	1889	1894	rVB5	8838	13397	0.72%	0.074%
34	14.310	1919	1925	1930	rBV5	18103	42370	2.28%	0.233%

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

35	14.451	1946	1949	1954	rVB6	4938	6107	0.33%	0.034%
36	14.875	2017	2021	2024	rVV4	3276	5224	0.28%	0.029%
37	14.934	2027	2031	2033	rVB4	3256	3618	0.19%	0.020%
38	15.039	2043	2049	2056	rVV	293210	424497	22.87%	2.339%
39	15.098	2056	2059	2065	rVB2	8188	12170	0.66%	0.067%
40	15.204	2073	2077	2089	rBV2	39383	76636	4.13%	0.422%
41	15.398	2107	2110	2114	rBV2	2168	3436	0.19%	0.019%
42	15.645	2148	2152	2157	rVB4	6629	8866	0.48%	0.049%
43	15.786	2172	2176	2177	rBV2	4207	4090	0.22%	0.023%
44	16.157	2235	2239	2241	rBV3	2897	4292	0.23%	0.024%
45	16.586	2307	2312	2315	rBV3	7700	11010	0.59%	0.061%
46	16.786	2338	2346	2351	rBV	1172139	1611913	86.85%	8.882%
47	16.828	2351	2353	2358	rVV	124925	158889	8.56%	0.876%
48	16.892	2358	2364	2374	rVV2	275555	444637	23.96%	2.450%
49	17.216	2413	2419	2426	rBV5	8973	17293	0.93%	0.095%
50	17.322	2434	2437	2441	rVB4	4710	5647	0.30%	0.031%
51	17.669	2493	2496	2502	rBV4	11300	17341	0.93%	0.096%
52	17.722	2502	2505	2509	rVV2	19152	25503	1.37%	0.141%
53	17.775	2509	2514	2524	rVV3	18036	40813	2.20%	0.225%
54	17.863	2524	2529	2533	rVV6	20445	30990	1.67%	0.171%
55	17.898	2533	2535	2541	rVB4	9607	15016	0.81%	0.083%
56	18.022	2550	2556	2559	rBV5	4570	8566	0.46%	0.047%
57	18.180	2580	2583	2588	rVB3	7084	10169	0.55%	0.056%
58	18.422	2620	2624	2628	rBV6	4308	6384	0.34%	0.035%
59	18.486	2630	2635	2637	rBV3	5520	6366	0.34%	0.035%
60	18.516	2637	2640	2644	rVV6	3770	5229	0.28%	0.029%
61	18.604	2649	2655	2660	rBV7	8754	15685	0.85%	0.086%
62	18.663	2661	2665	2669	rBV5	8347	11197	0.60%	0.062%
63	18.745	2676	2679	2683	rVB5	7370	11467	0.62%	0.063%
64	18.827	2689	2693	2694	rBV4	6310	9077	0.49%	0.050%
65	18.863	2694	2699	2708	rVV	225371	311451	16.78%	1.716%
66	18.951	2710	2714	2715	rBV4	4568	5254	0.28%	0.029%
67	19.192	2751	2755	2759	rBV	413401	629911	33.94%	3.471%
68	19.304	2772	2774	2776	rBV3	8163	8169	0.44%	0.045%
69	19.357	2781	2783	2786	rBV4	4831	5851	0.32%	0.032%
70	19.422	2790	2794	2798	rBV4	13301	19004	1.02%	0.105%
71	19.569	2812	2819	2826	rBV10	17006	48662	2.62%	0.268%

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72	19.692	2838	2840	2842	rBV2	8401	8664	0.47%	0.048%
73	19.739	2844	2848	2853	rVB4	24362	35481	1.91%	0.196%
74	19.792	2853	2857	2860	rVB5	13933	17226	0.93%	0.095%
75	19.839	2861	2865	2868	rBV4	16541	23077	1.24%	0.127%
76	19.898	2871	2875	2879	rBV5	21267	33153	1.79%	0.183%
77	20.010	2892	2894	2895	rBV2	7204	5122	0.28%	0.028%
78	20.039	2895	2899	2903	rVB4	9876	15175	0.82%	0.084%
79	20.292	2939	2942	2945	rBV3	19499	24459	1.32%	0.135%
80	20.657	3000	3004	3006	rBV3	19449	25022	1.35%	0.138%
81	20.698	3006	3011	3015	rBV2	110750	196186	10.57%	1.081%
82	20.751	3016	3020	3025	rVV2	245331	363312	19.58%	2.002%
83	20.804	3025	3029	3038	rVV	329539	436423	23.51%	2.405%
84	20.874	3038	3041	3044	rVV3	22130	31603	1.70%	0.174%
85	20.951	3051	3054	3057	rBV	49929	62438	3.36%	0.344%
86	21.004	3057	3063	3067	rVV	1358496	1855939	100.00%	10.227%
87	21.039	3067	3069	3076	rVB	131156	169316	9.12%	0.933%
88	21.192	3092	3095	3099	rBV2	75120	73785	3.98%	0.407%
89	21.610	3163	3166	3173	rBV	139264	165330	8.91%	0.911%
90	22.045	3236	3240	3248	rVB	166444	215472	11.61%	1.187%
91	22.251	3272	3275	3281	rVB3	118870	150763	8.12%	0.831%
92	22.510	3312	3319	3322	rBV2	329738	583988	31.47%	3.218%
93	22.545	3323	3325	3332	rVB3	190407	238799	12.87%	1.316%
94	22.974	3393	3398	3402	rBV	279861	441508	23.79%	2.433%
95	23.021	3402	3406	3412	rVB2	270057	453485	24.43%	2.499%
96	23.104	3414	3420	3433	rVB	1004088	1693968	91.27%	9.334%
97	23.610	3501	3506	3512	rVB	363951	584440	31.49%	3.220%
98	23.680	3514	3518	3523	rVV5	64935	114150	6.15%	0.629%
99	24.274	3615	3619	3630	rVB	172414	350824	18.90%	1.933%
100	25.057	3747	3752	3760	rVB2	174658	385525	20.77%	2.124%

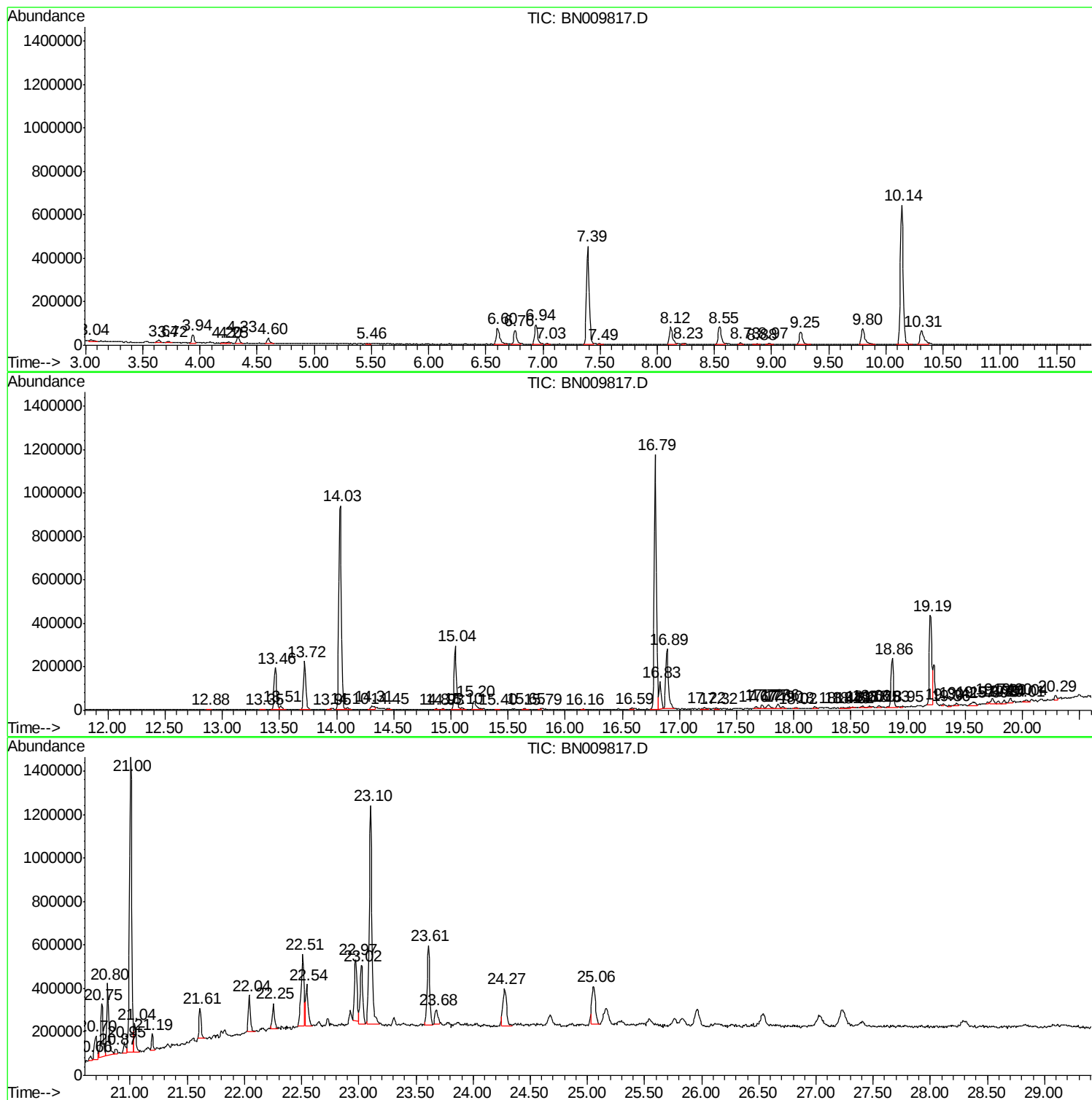
Sum of corrected areas: 18148071

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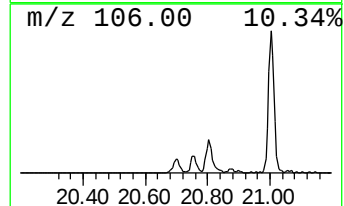
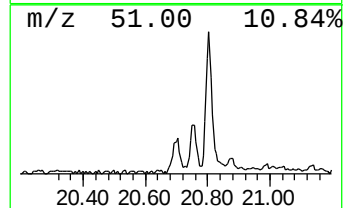
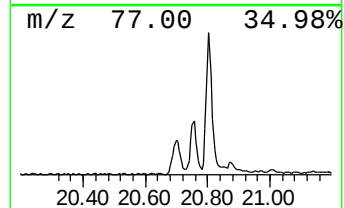
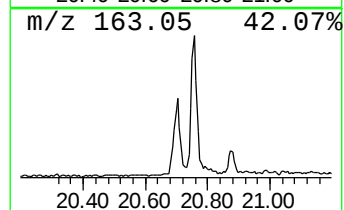
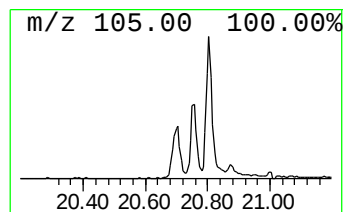
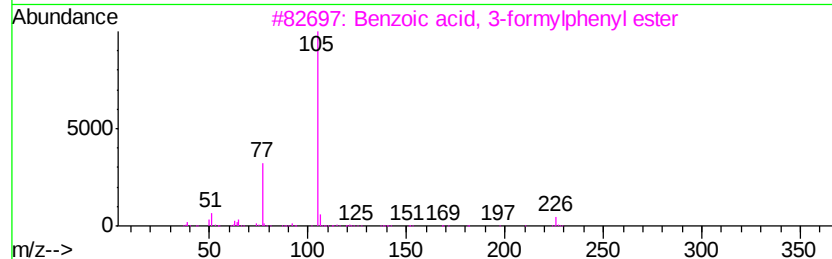
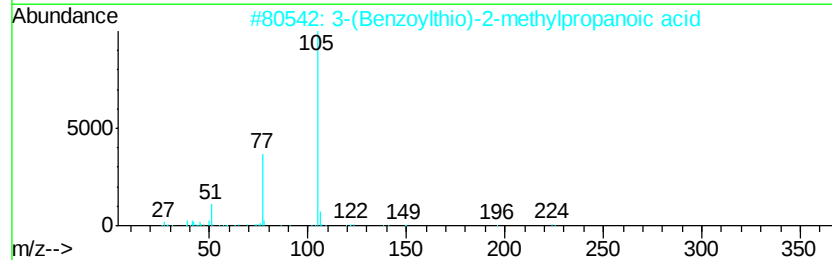
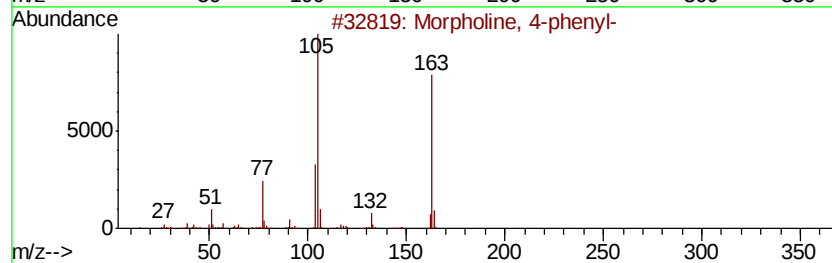
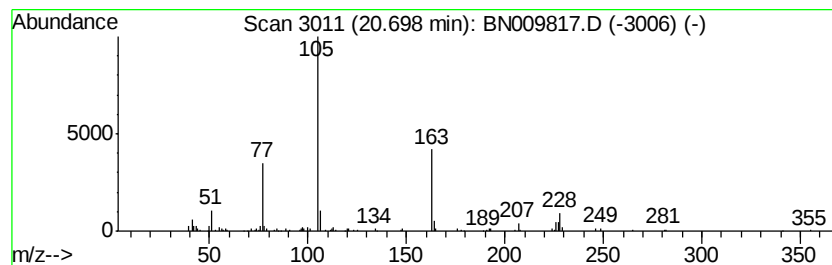
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN021220MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 unknown-01 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.70	2.11 ng/ul	196186	Chrysene-d12	21.00

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Morpholine, 4-phenyl-	163	C10H13NO	000092-53-5	47
2		3-(Benzoylthio)-2-methylpropanoi...	224	C11H12O3S	067714-34-5	41
3		Benzoic acid, 3-formylphenyl ester	226	C14H10O3	033696-06-9	38
4		2-Methoxyphenyl benzoate	228	C14H12O3	000531-37-3	38
5		2-Methoxyphenyl benzoate	228	C14H12O3	000531-37-3	38



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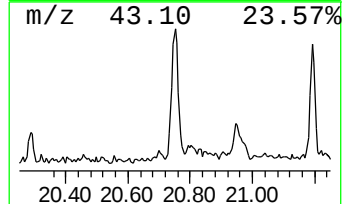
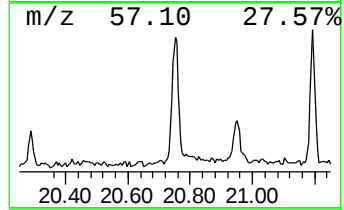
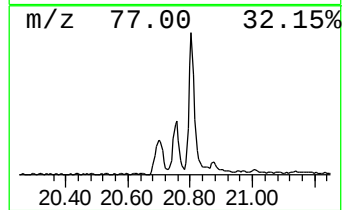
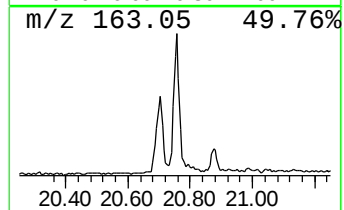
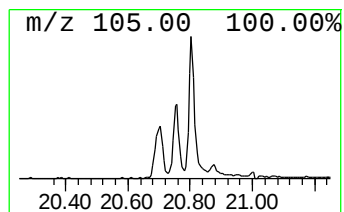
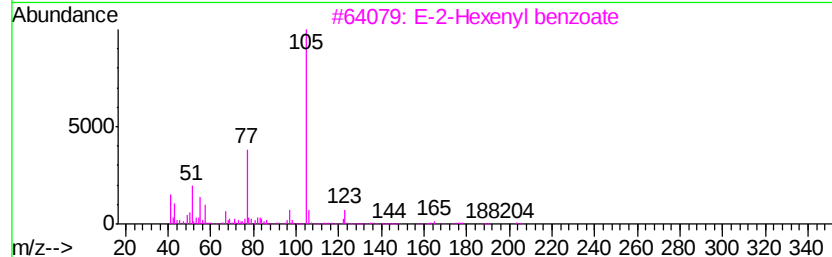
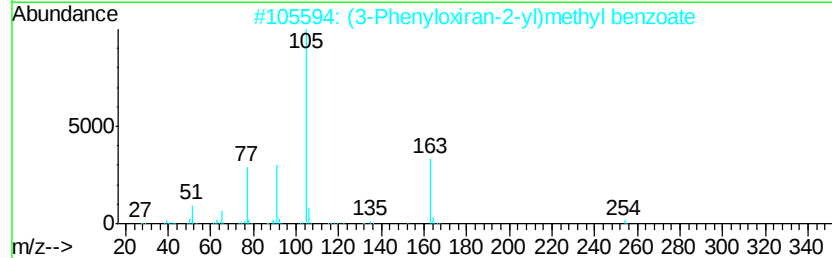
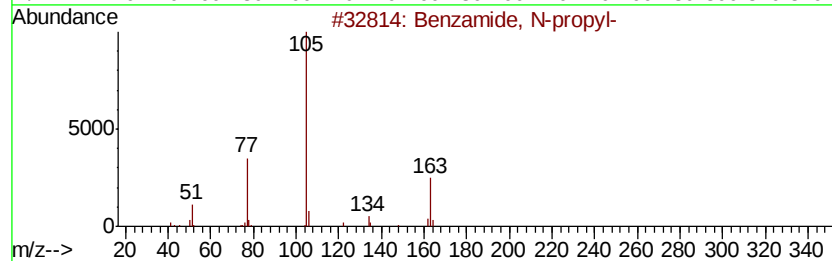
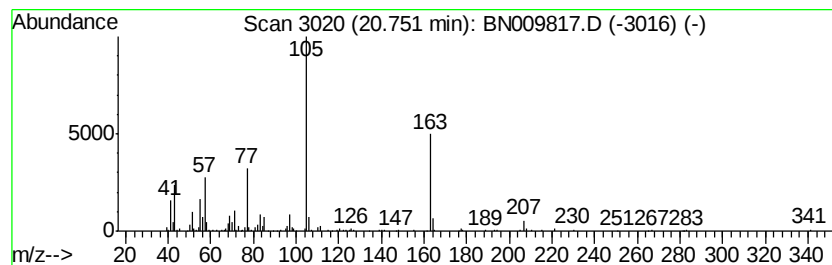
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN021220MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 Benzamide, N-propyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.75	3.92 ng/ul	363312	Chrysene-d12	21.00
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Benzamide, N-propyl-	163 C10H13NO	010546-70-0 64
2		(3-Phenyloxiran-2-yl)methyl benz...	254 C16H14O3	1000366-96-1 47
3		E-2-Hexenyl benzoate	204 C13H16O2	076841-70-8 35
4		2-Bromoisobutyrophenone	226 C10H11BrO	010409-54-8 22
5		Benzoic acid 5-methyl-2-phenyl-2...	278 C17H14N2O2	056159-67-2 22



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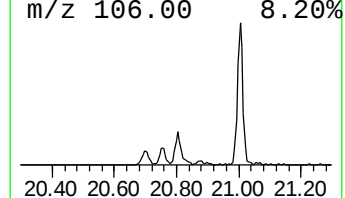
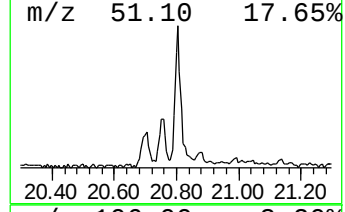
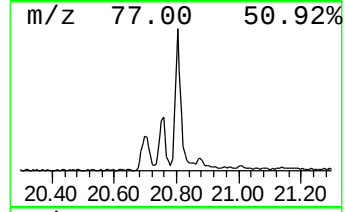
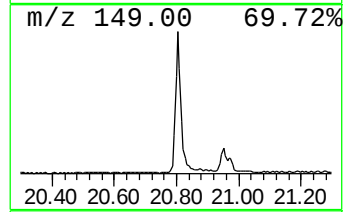
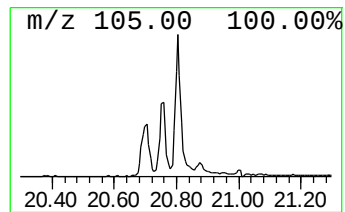
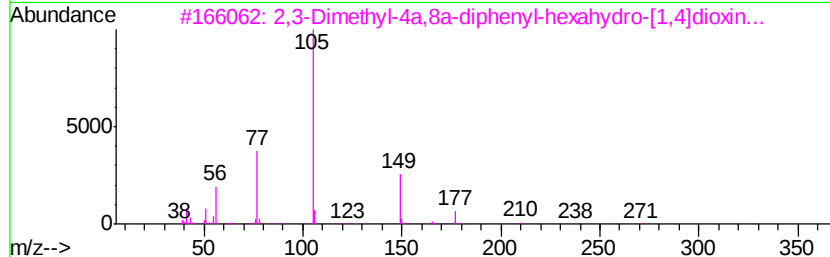
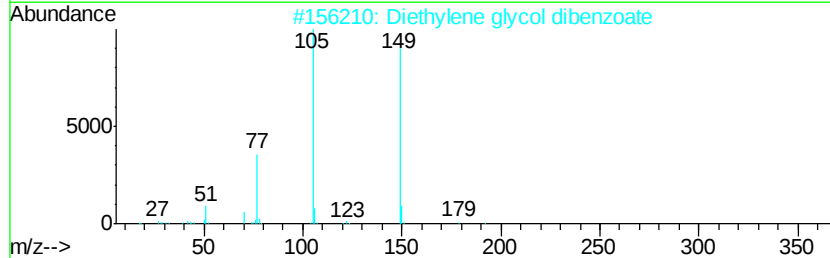
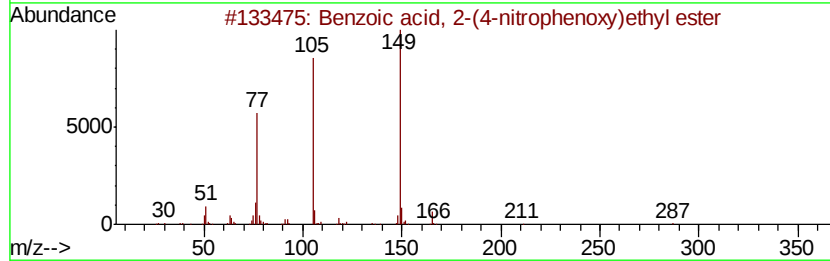
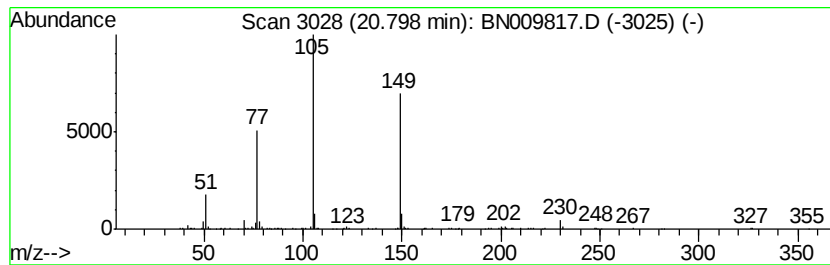
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Quant Title : SVOA CALIBRATION

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

Peak Number 3 Benzoic acid, 2-(4-nitrophe... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.		
20.80	4.70 ng/ul	436423	Chrysene-d12	21.00		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzoic acid, 2-(4-nitrophenoxyv)...	287	C15H13N05	1000368-92-7	72	
2	Diethylene glycol dibenzoate	314	C18H1805	000120-55-8	64	
3	2,3-Dimethyl-4a,8a-diphenyl-hexa...	326	C20H2204	1000297-07-0	43	
4	2-Methyl-4-hydroxybenzoxazole	149	C8H7N02	051110-60-2	43	
5	Vinyl benzoate	148	C9H802	000769-78-8	38	



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN022220\
Data File : BN009817.D
Acq On : 21 Feb 2020 20:55
Operator : CG/JU
Sample : L1535-03
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
CB6Q1

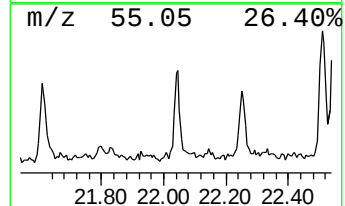
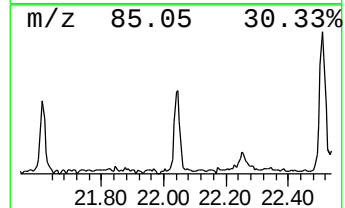
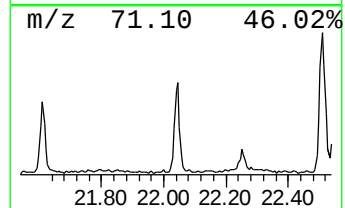
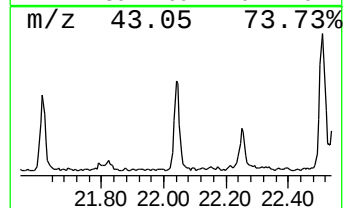
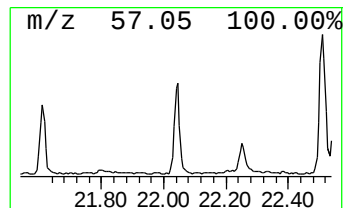
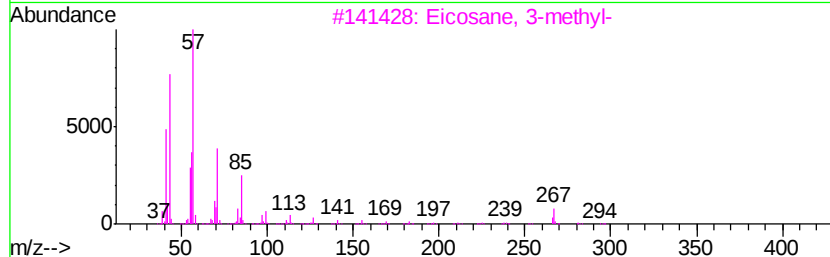
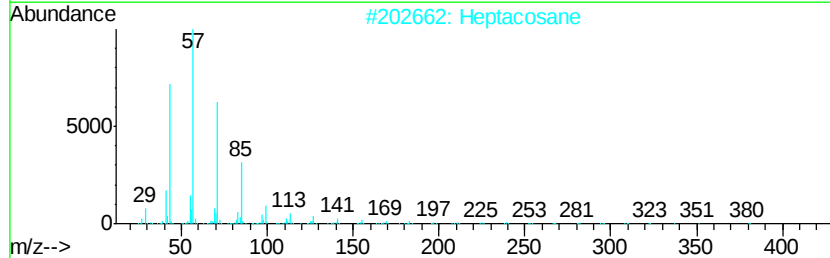
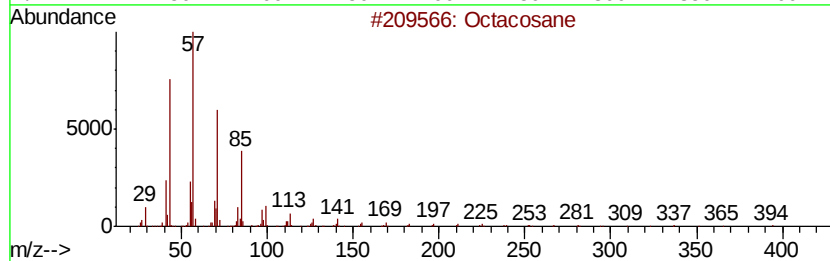
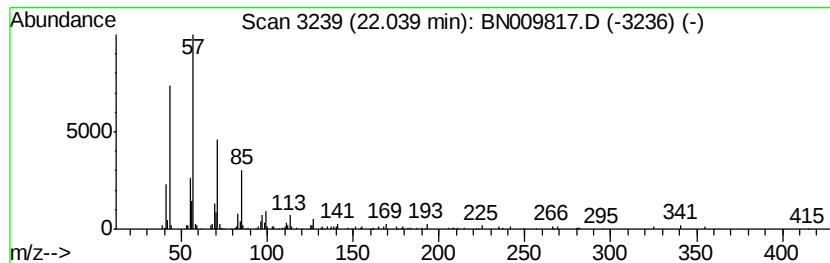
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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.04	2.32 ng/ul	215472	Chrysene-d12	21.00

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octacosane	394	C28H58	000630-02-4	90
2		Heptacosane	380	C27H56	000593-49-7	87
3		Eicosane, 3-methyl-	296	C21H44	006418-46-8	86
4		Heptadecane, 9-hexyl-	324	C23H48	055124-79-3	80
5		Heptacosane	380	C27H56	000593-49-7	80



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN022220\
Data File : BN009817.D
Acq On : 21 Feb 2020 20:55
Operator : CG/JU
Sample : L1535-03
Misc :
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Instrument :
BNA_N
ClientSampleId :
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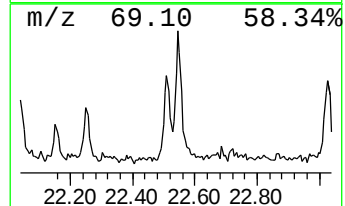
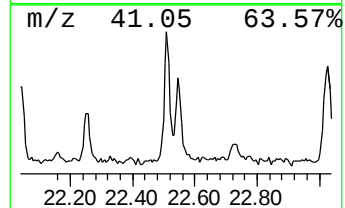
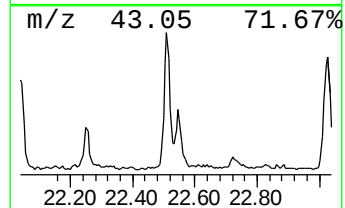
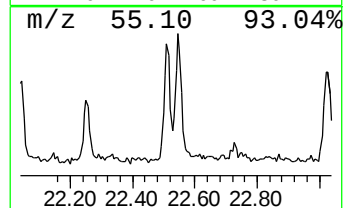
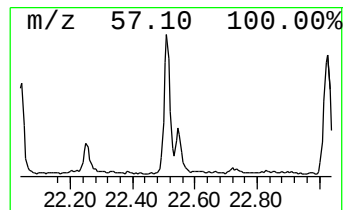
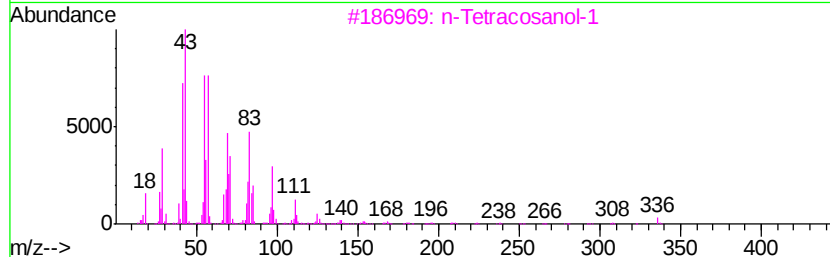
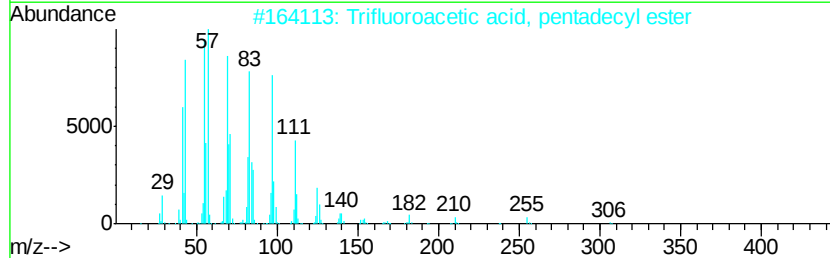
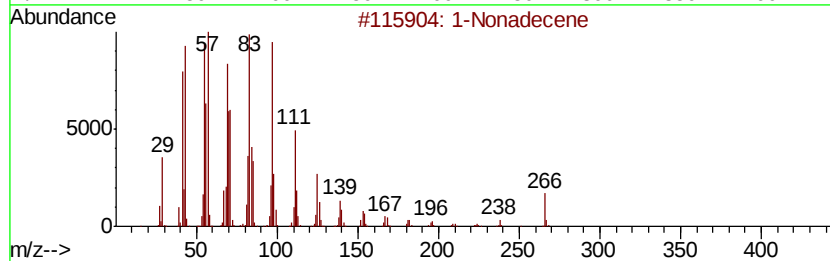
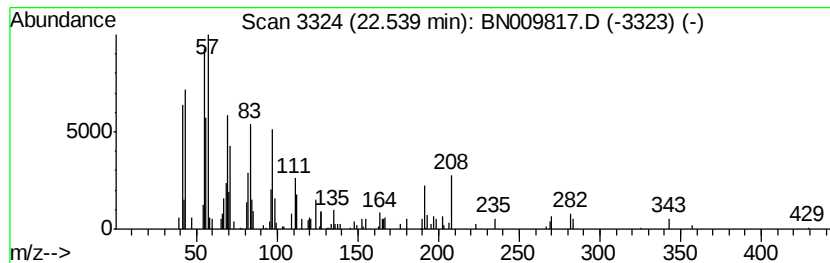
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 (DEL) Alkane: Straight-Chai... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.54	2.82 ng/ul	238799	Perylene-d12	23.10

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1-Nonadecene	266	C19H38	018435-45-5	83
2		Trifluoroacetic acid, pentadecyl...	324	C17H31F3O2	959010-23-2	72
3		n-Tetracosanol-1	354	C24H50O	000506-51-4	62
4		Heptafluorobutyric acid, pentade...	424	C19H31F7O2	959261-23-5	58
5		1-Hexacosanol	382	C26H54O	000506-52-5	58



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN022220\
Data File : BN009817.D
Acq On : 21 Feb 2020 20:55
Operator : CG/JU
Sample : L1535-03
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
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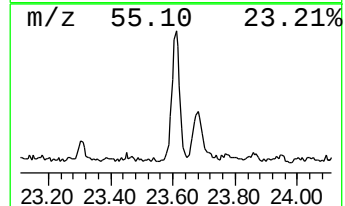
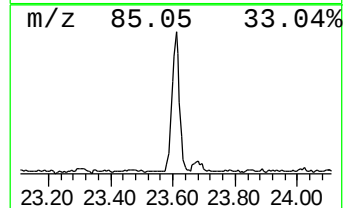
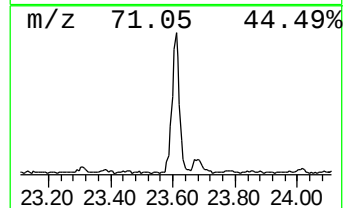
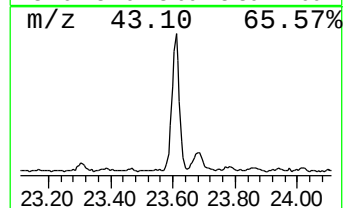
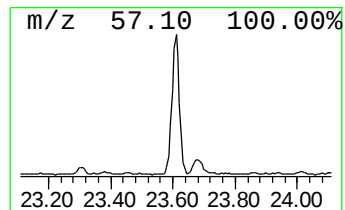
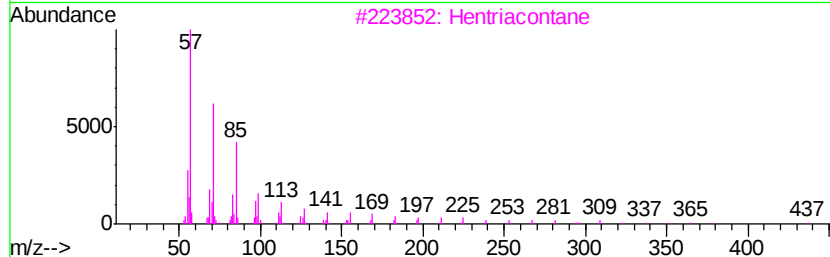
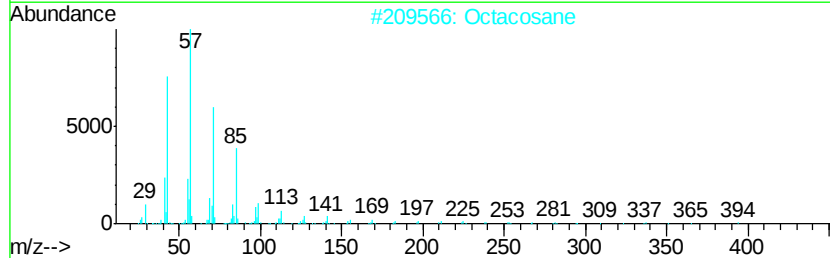
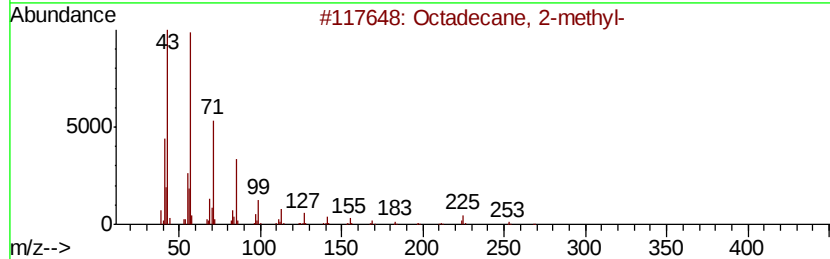
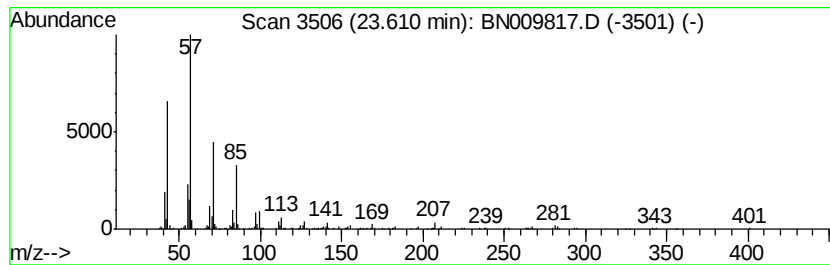
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 (DEL) Alkane: Straight-Chai... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.61	6.90 ng/ul	584440	Perylene-d12	23.10

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octadecane, 2-methyl-	268	C19H40	001560-88-9	94
2			Octacosane	394	C28H58	000630-02-4	87
3			Hentriacontane	437	C31H64	000630-04-6	87
4			Hentriacontane	437	C31H64	000630-04-6	87
5			Nonadecane, 9-methyl-	282	C20H42	013287-24-6	83



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN022220\
Data File : BN009817.D
Acq On : 21 Feb 2020 20:55
Operator : CG/JU
Sample : L1535-03
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
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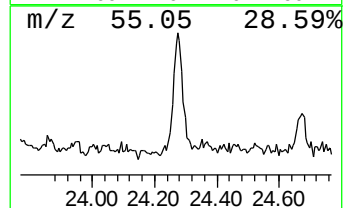
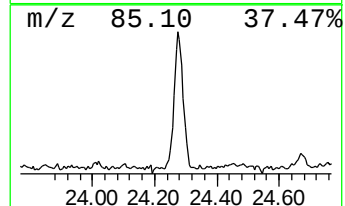
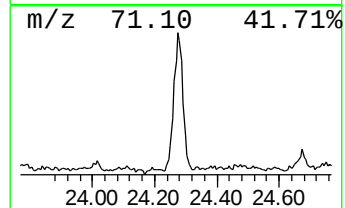
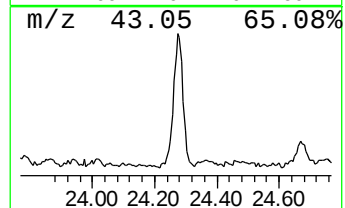
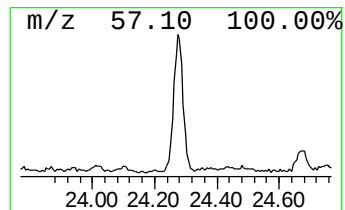
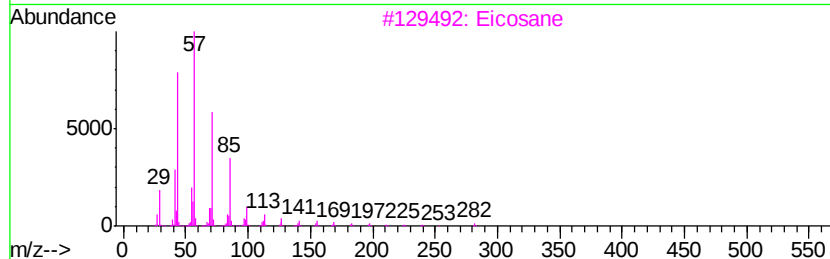
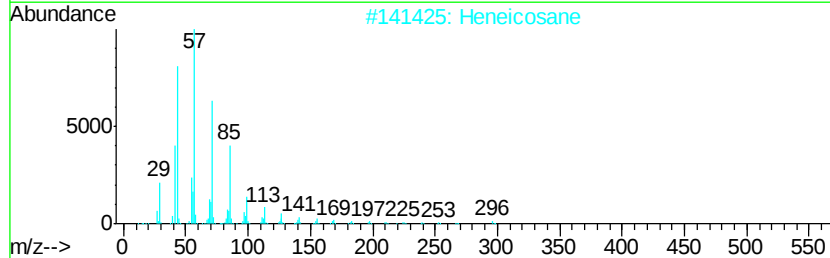
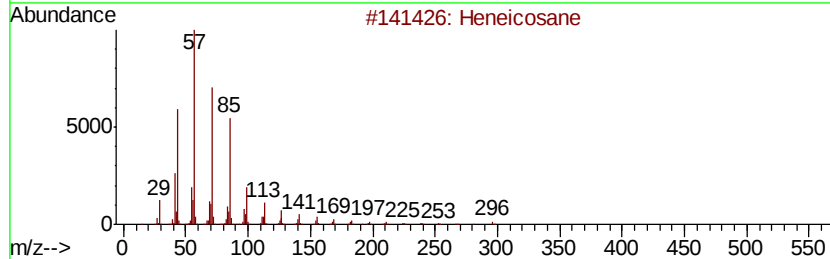
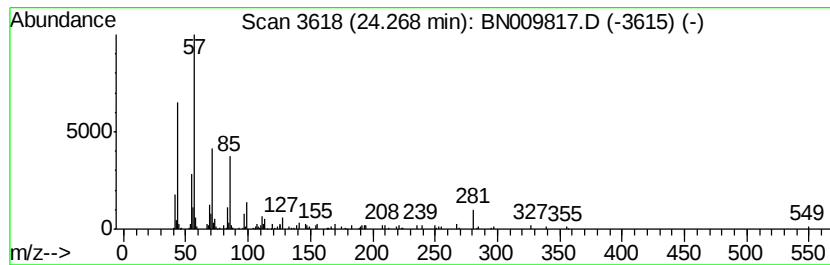
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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.27	4.14 ng/ul	350824	Perylene-d12	23.10

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heneicosane	296	C21H44	000629-94-7	97
2		Heneicosane	296	C21H44	000629-94-7	95
3		Eicosane	282	C20H42	000112-95-8	90
4		Eicosane	282	C20H42	000112-95-8	86
5		Eicosane	282	C20H42	000112-95-8	86



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN022220\
Data File : BN009817.D
Acq On : 21 Feb 2020 20:55
Operator : CG/JU
Sample : L1535-03
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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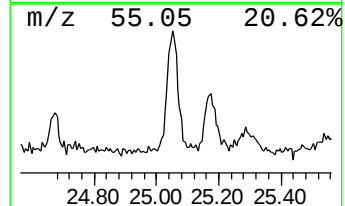
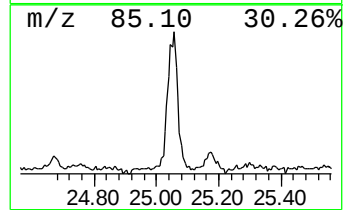
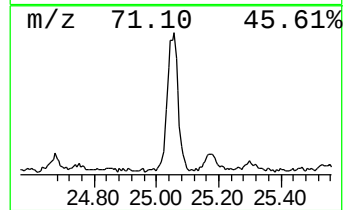
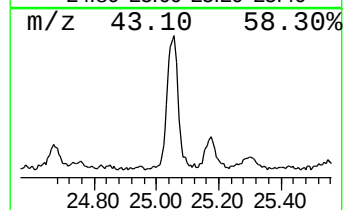
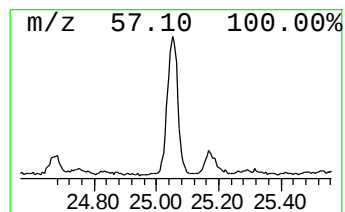
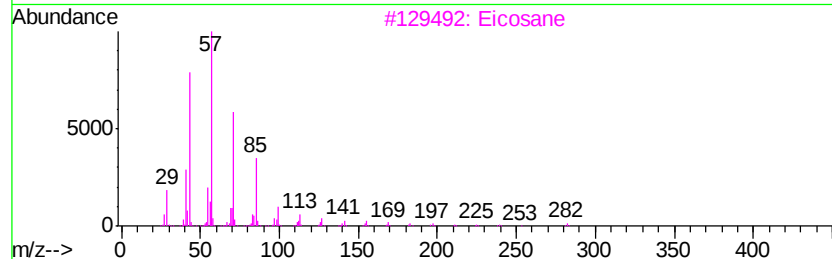
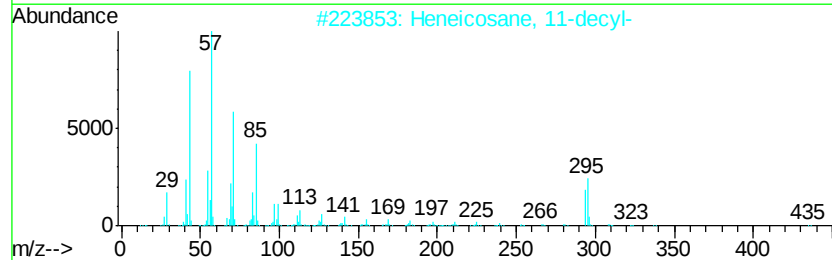
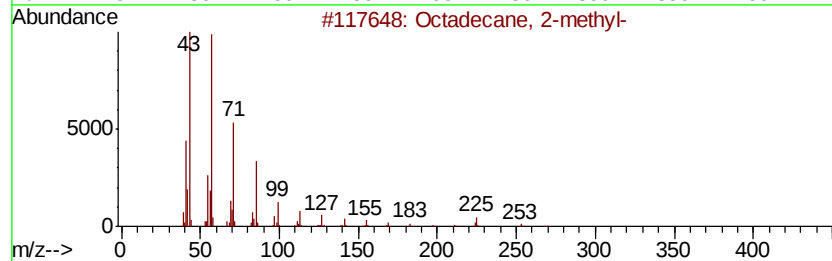
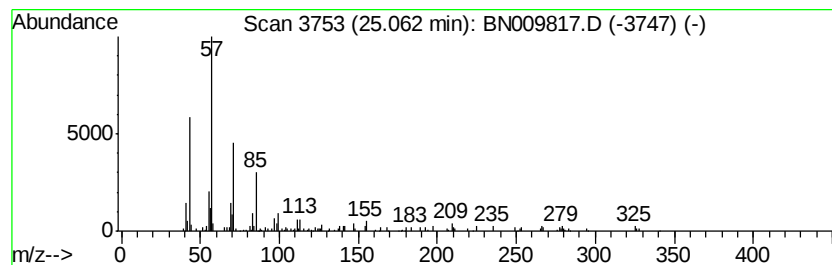
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Quant Title : SVOA CALIBRATION

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.
25.06	4.55 ng/ul	385525	Perylene-d12	23.10

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octadecane, 2-methyl-	268	C19H40	001560-88-9	86
2		Heneicosane, 11-decyl-	437	C31H64	055320-06-4	74
3		Eicosane	282	C20H42	000112-95-8	74
4		Undecane, 3,8-dimethyl-	184	C13H28	017301-30-3	72
5		Sulfurous acid, butyl dodecyl ester	306	C16H34O3S	1000309-17-9	72



Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN022220\
Data File : BN009817.D
Acq On : 21 Feb 2020 20:55
Operator : CG/JU
Sample : L1535-03
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
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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
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Benzamide, N-propyl-	20.75	3.9	ng/ul	363312	5	21.00	1855940	20.0
Benzoic acid, 2-(...	20.80	4.7	ng/ul	436423	5	21.00	1855940	20.0
(DEL) Alkane: Str...	22.04	2.3	ng/ul	215472	5	21.00	1855940	20.0
(DEL) Alkane: Str...	22.54	2.8	ng/ul	238799	6	23.10	1693970	20.0
(DEL) Alkane: Str...	23.61	6.9	ng/ul	584440	6	23.10	1693970	20.0
(DEL) Alkane: Str...	24.27	4.1	ng/ul	350824	6	23.10	1693970	20.0
(DEL) Alkane: Str...	25.06	4.5	ng/ul	385525	6	23.10	1693970	20.0