

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN022220\
 Data File : BN009831.D
 Acq On : 22 Feb 2020 05:51
 Operator : CG/JU
 Sample : L1535-14
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 CB6R2

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	3.711	120	123	129	rVB	47124	63414	2.68%	0.224%
2	3.934	157	161	169	rVB2	90573	121801	5.15%	0.430%
3	4.069	182	184	191	rVB5	13455	17272	0.73%	0.061%
4	4.328	223	228	236	rBV	72284	108938	4.60%	0.385%
5	4.587	268	272	276	rBV	36518	55662	2.35%	0.197%
6	5.464	417	421	426	rVB5	10171	14738	0.62%	0.052%
7	5.522	426	431	435	rBV6	9403	18273	0.77%	0.065%
8	6.099	526	529	534	rVB2	15773	21803	0.92%	0.077%
9	6.599	606	614	628	rVB	215583	399267	16.87%	1.411%
10	6.752	634	640	649	rBV	160558	257154	10.87%	0.909%
11	6.934	665	671	682	rBV	230778	373370	15.78%	1.319%
12	7.028	685	687	695	rVB5	10265	17189	0.73%	0.061%
13	7.111	697	701	707	rBV6	7394	11908	0.50%	0.042%
14	7.387	742	748	758	rBV	461134	757998	32.03%	2.678%
15	7.481	761	764	768	rVB2	6931	10332	0.44%	0.037%
16	8.105	865	870	882	rVV	236332	427851	18.08%	1.512%
17	8.216	886	889	900	rVB3	10147	23895	1.01%	0.084%
18	8.540	937	944	953	rVV	198777	369530	15.61%	1.306%
19	8.610	953	956	961	rVB5	9333	14833	0.63%	0.052%
20	8.710	968	973	980	rBV	52321	89092	3.76%	0.315%
21	8.975	1013	1018	1024	rVB	52052	74733	3.16%	0.264%
22	9.246	1059	1064	1073	rBV	167156	296708	12.54%	1.048%
23	9.563	1114	1118	1123	rBV4	6832	9974	0.42%	0.035%
24	9.781	1149	1155	1170	rBV	222645	431383	18.23%	1.524%
25	10.140	1208	1216	1222	rBV	659136	1099668	46.46%	3.886%
26	10.187	1222	1224	1230	rVB2	23603	28490	1.20%	0.101%
27	10.299	1234	1243	1264	rBV	170053	408416	17.26%	1.443%
28	11.599	1457	1464	1471	rVB5	11323	22689	0.96%	0.080%
29	11.698	1474	1481	1486	rBV4	12818	25889	1.09%	0.091%
30	11.740	1486	1488	1493	rVV5	7072	10280	0.43%	0.036%
31	11.822	1495	1502	1507	rVV4	10644	19307	0.82%	0.068%
32	12.040	1535	1539	1544	rBV5	7270	11341	0.48%	0.040%
33	12.869	1676	1680	1685	rVB2	16886	22195	0.94%	0.078%
34	13.351	1756	1762	1767	rBV7	11230	17216	0.73%	0.061%

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

35	13.457	1774	1780	1785	rBV	450747	639449	27.02%	2.259%
36	13.504	1785	1788	1793	rVB	35803	52621	2.22%	0.186%
37	13.593	1798	1803	1807	rVB5	5833	10228	0.43%	0.036%
38	13.716	1818	1824	1834	rBV	516052	789998	33.38%	2.791%
39	13.963	1863	1866	1871	rVB2	14207	17763	0.75%	0.063%
40	14.028	1871	1877	1885	rBV2	942816	1404727	59.35%	4.963%
41	14.092	1885	1888	1894	rVB	20830	29926	1.26%	0.106%
42	14.287	1914	1921	1932	rBV	83320	216348	9.14%	0.764%
43	14.445	1944	1948	1954	rVB2	15307	24848	1.05%	0.088%
44	14.657	1981	1984	1991	rVB7	5277	9474	0.40%	0.033%
45	14.922	2026	2029	2034	rVB4	14707	21692	0.92%	0.077%
46	15.034	2042	2048	2054	rBV	678413	963407	40.70%	3.404%
47	15.092	2054	2058	2067	rVB	27707	46130	1.95%	0.163%
48	15.192	2071	2075	2084	rBV2	146116	265030	11.20%	0.936%
49	15.334	2097	2099	2104	rVB5	7110	9307	0.39%	0.033%
50	15.392	2104	2109	2113	rBV5	10389	16002	0.68%	0.057%
51	15.539	2127	2134	2141	rVV6	16673	34009	1.44%	0.120%
52	15.639	2144	2151	2155	rVV5	18740	32049	1.35%	0.113%
53	15.787	2169	2176	2180	rBV6	19739	43764	1.85%	0.155%
54	15.957	2201	2205	2209	rBV5	6830	11028	0.47%	0.039%
55	16.151	2235	2238	2243	rVB6	9411	13600	0.57%	0.048%
56	16.334	2266	2269	2273	rBV5	6851	11491	0.49%	0.041%
57	16.457	2286	2290	2296	rVB4	17447	29842	1.26%	0.105%
58	16.539	2298	2304	2307	rBV7	7841	17315	0.73%	0.061%
59	16.586	2307	2312	2314	rBV3	20146	32114	1.36%	0.113%
60	16.786	2340	2346	2350	rBV	1192627	1618385	68.38%	5.718%
61	16.828	2350	2353	2358	rVB	383464	491470	20.77%	1.737%
62	16.886	2358	2363	2368	rBV2	683968	983988	41.57%	3.477%
63	17.204	2413	2417	2422	rBV2	37851	55554	2.35%	0.196%
64	17.322	2433	2437	2442	rVB5	18006	20423	0.86%	0.072%
65	17.498	2463	2467	2475	rVB10	9009	19397	0.82%	0.069%
66	17.663	2491	2495	2500	rVV	51373	72758	3.07%	0.257%
67	17.716	2500	2504	2509	rVB2	82598	113885	4.81%	0.402%
68	17.857	2523	2528	2532	rBV3	81010	131263	5.55%	0.464%
69	17.898	2532	2535	2540	rVB	39112	54740	2.31%	0.193%
70	18.145	2575	2577	2578	rBV2	13391	8971	0.38%	0.032%
71	18.175	2579	2582	2587	rVV2	45730	66900	2.83%	0.236%

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72	18.228	2587	2591	2595	rVB4	21191	31662	1.34%	0.112%
73	18.475	2630	2633	2637	rBV4	19121	23946	1.01%	0.085%
74	18.516	2638	2640	2643	rVB4	17632	19203	0.81%	0.068%
75	18.598	2649	2654	2659	rBV2	51182	82394	3.48%	0.291%
76	18.657	2659	2664	2668	rBV5	61756	91568	3.87%	0.324%
77	18.739	2674	2678	2684	rBV5	36739	66334	2.80%	0.234%
78	18.857	2693	2698	2707	rVB	734543	976003	41.24%	3.449%
79	19.016	2722	2725	2730	rBV3	24192	31728	1.34%	0.112%
80	19.192	2750	2755	2758	rBV	964141	1400399	59.17%	4.948%
81	19.222	2758	2760	2769	rVB	682348	832098	35.16%	2.940%
82	19.416	2791	2793	2799	rVB	41563	46987	1.99%	0.166%
83	19.563	2813	2818	2823	rBV5	53602	119068	5.03%	0.421%
84	19.698	2838	2841	2842	rBV3	39788	43541	1.84%	0.154%
85	19.786	2853	2856	2861	rVB	63036	67930	2.87%	0.240%
86	19.833	2861	2864	2868	rBV	62322	78067	3.30%	0.276%
87	19.892	2870	2874	2878	rVV2	68336	93409	3.95%	0.330%
88	20.575	2987	2990	2995	rVB	115472	137651	5.82%	0.486%
89	20.692	3007	3010	3014	rBV2	112206	181707	7.68%	0.642%
90	20.751	3016	3020	3024	rVV2	522212	768689	32.48%	2.716%
91	20.804	3025	3029	3037	rVB2	317044	444469	18.78%	1.570%
92	21.004	3057	3063	3067	rBV2	1520164	2366807	100.00%	8.363%
93	21.039	3067	3069	3075	rVB	365498	431493	18.23%	1.525%
94	21.192	3093	3095	3099	rVB	111092	106566	4.50%	0.377%
95	21.616	3164	3167	3177	rVB2	390843	608851	25.72%	2.151%
96	22.510	3312	3319	3322	rBV2	685078	1406570	59.43%	4.970%
97	22.974	3394	3398	3402	rBV	576149	823979	34.81%	2.911%
98	23.104	3415	3420	3425	rVB	923506	1486258	62.80%	5.251%
99	23.610	3502	3506	3512	rVB	526573	864327	36.52%	3.054%
100	23.680	3515	3518	3528	rVB	367830	637549	26.94%	2.253%

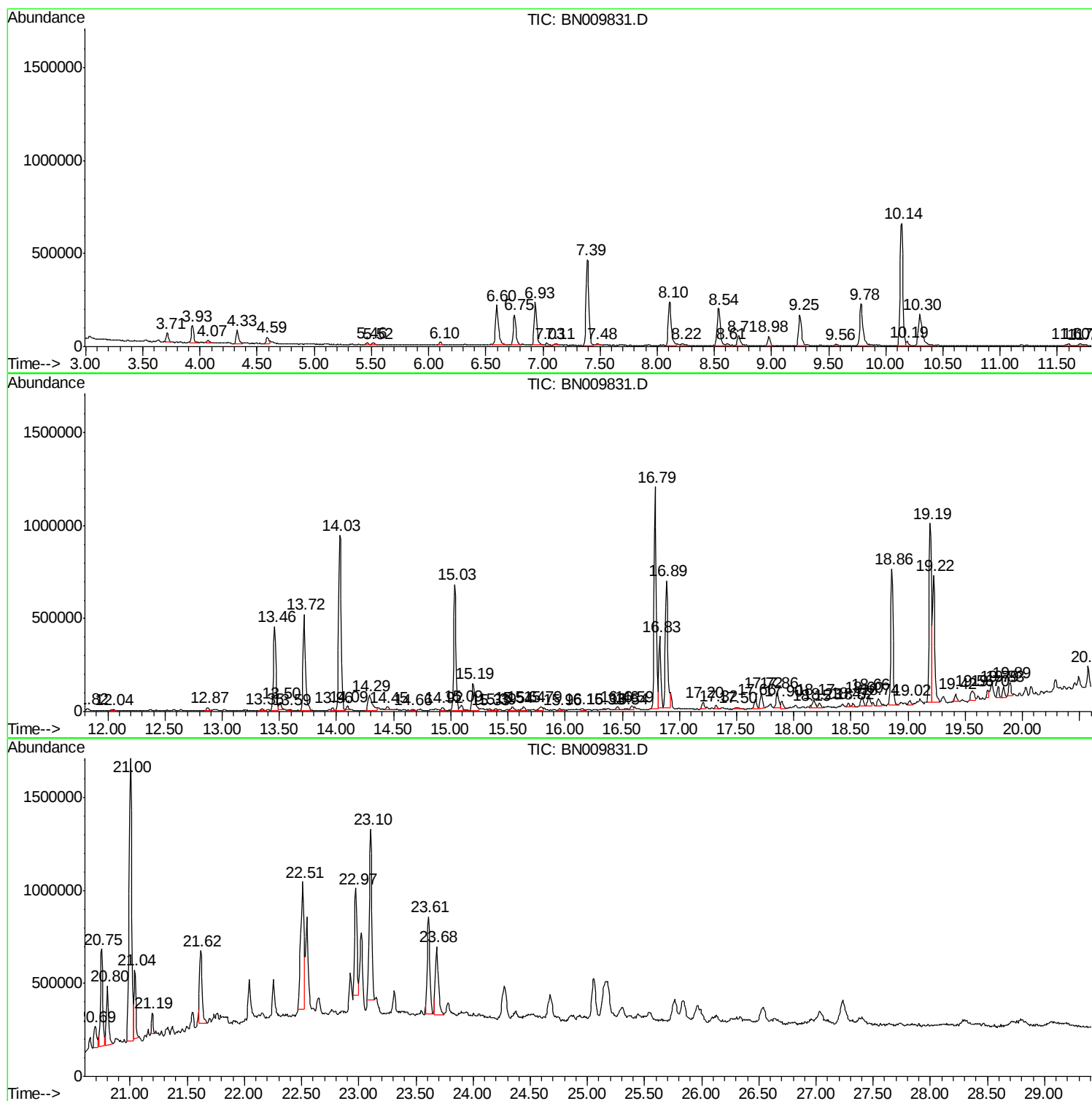
Sum of corrected areas: 28301758

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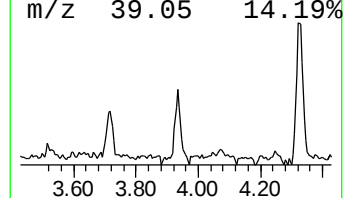
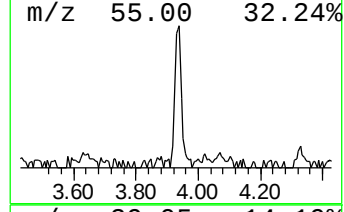
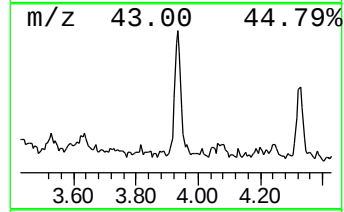
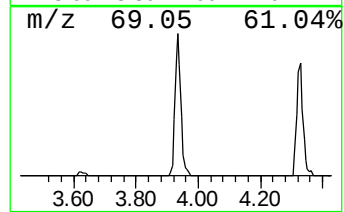
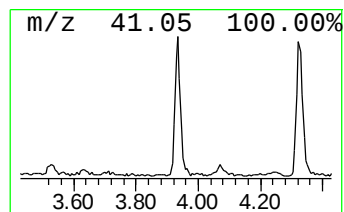
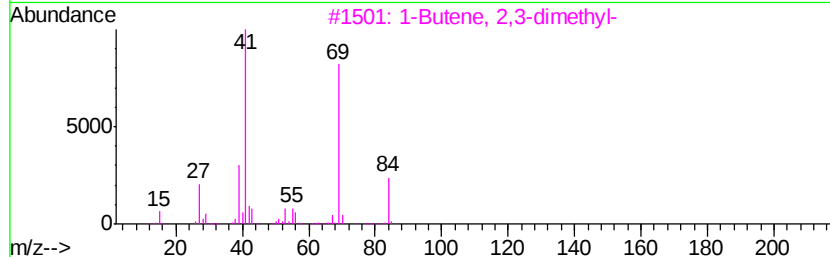
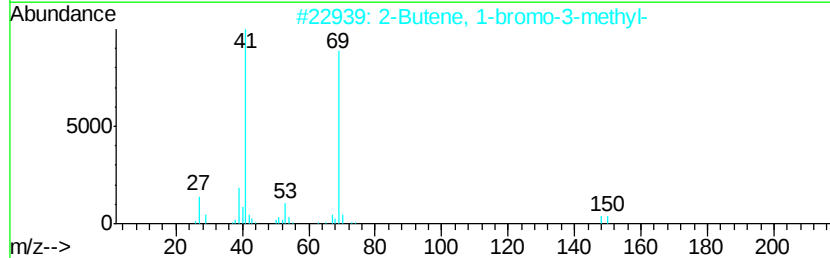
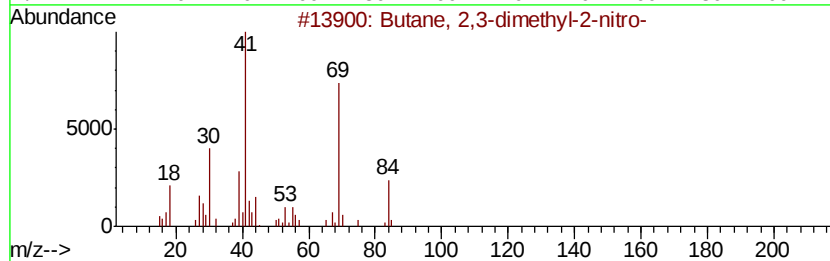
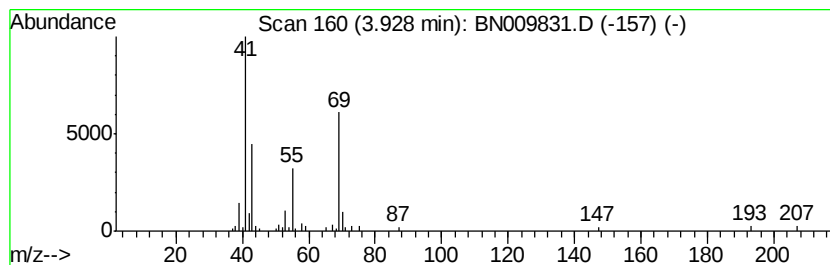
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN021220MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Butane, 2,3-dimethyl-2-nitro- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.93	3.21 ng/ul	121801	1,4-Dichlorobenzene-d4	7.39

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Butane, 2,3-dimethyl-2-nitro-	131	C6H13NO2	034075-28-0	53
2		2-Butene, 1-bromo-3-methyl-	148	C5H9Br	000870-63-3	47
3		1-Butene, 2,3-dimethyl-	84	C6H12	000563-78-0	42
4		2-Pentene, 2-methyl-	84	C6H12	000625-27-4	42
5		2-Pentene, 2-methyl-	84	C6H12	000625-27-4	42



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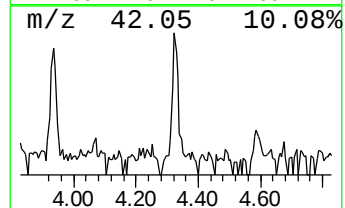
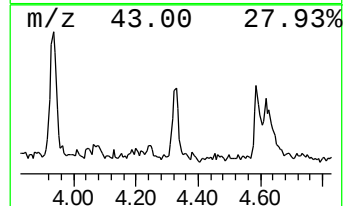
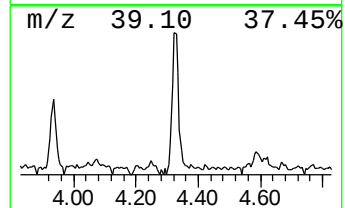
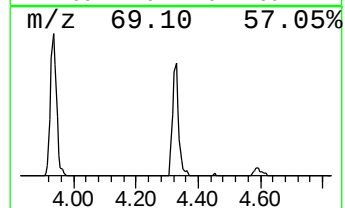
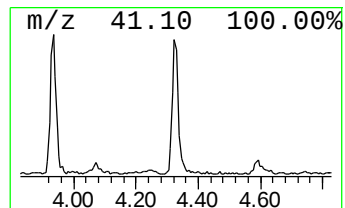
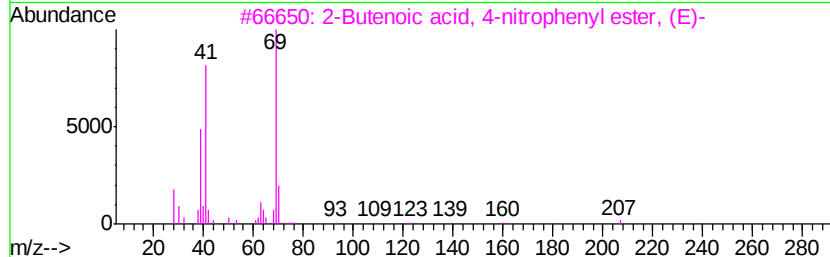
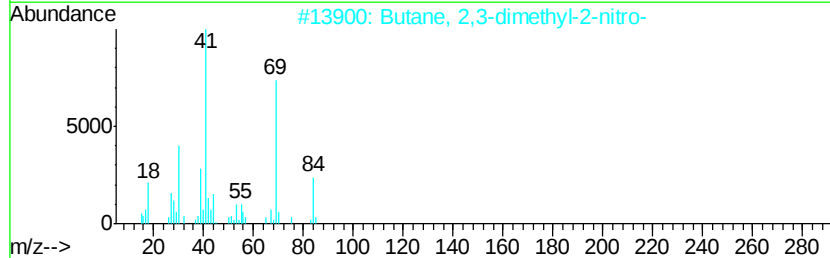
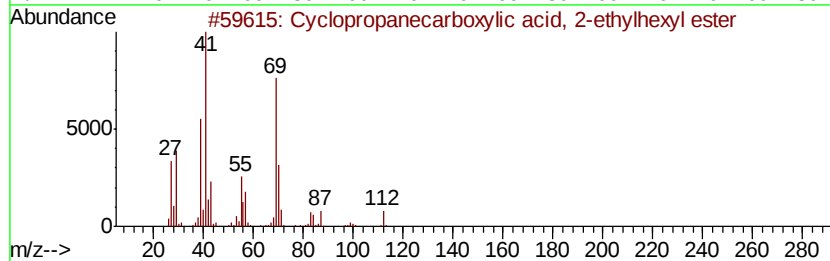
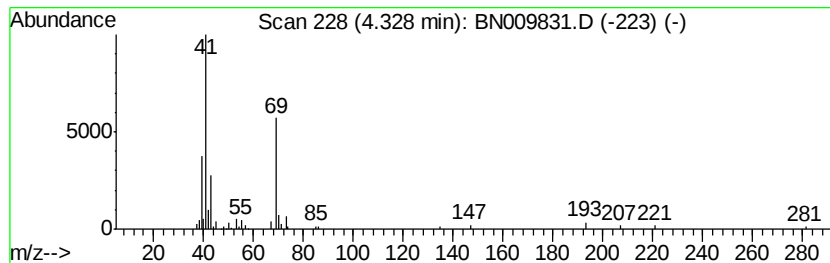
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN021220MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 unknown-01 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.33	2.87 ng/ul	108938	1,4-Dichlorobenzene-d4	7.39

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclopropanecarboxylic acid, 2-e...	198	C12H22O2	103604-31-5	28
2		Butane, 2,3-dimethyl-2-nitro-	131	C6H13NO2	034075-28-0	28
3		2-Butenoic acid, 4-nitrophenyl e...	207	C10H9NO4	014617-88-0	27
4		2-Butene, 1-bromo-3-methyl-	148	C5H9Br	000870-63-3	25
5		Cyclopentane, bromo-	148	C5H9Br	000137-43-9	17



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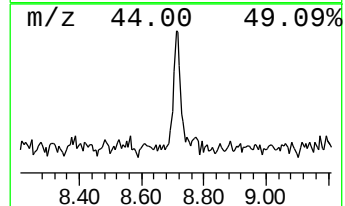
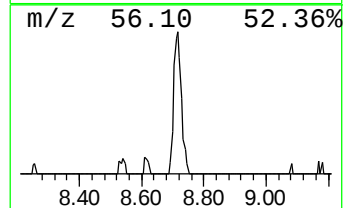
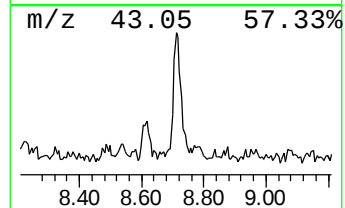
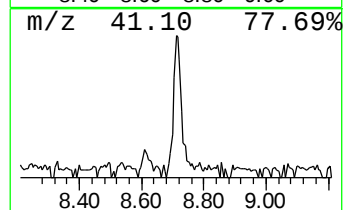
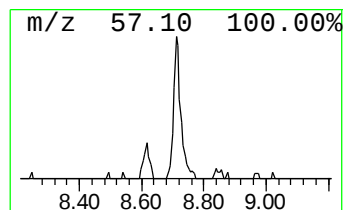
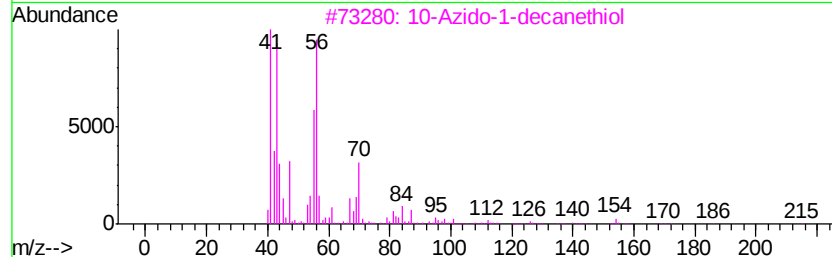
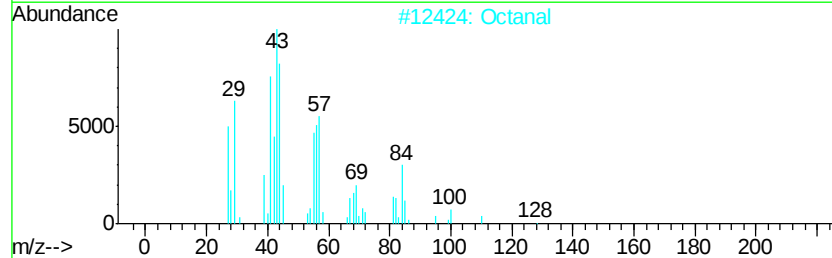
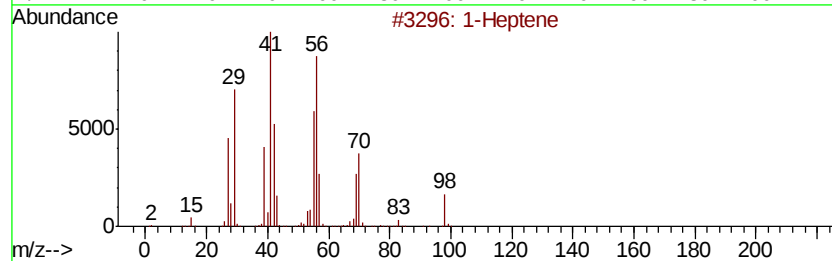
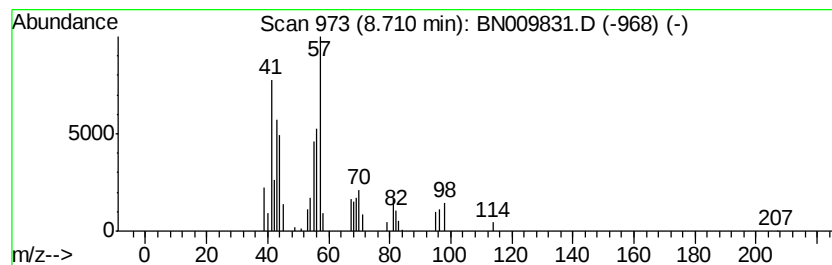
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.71	2.35 ng/ul	89092	1,4-Dichlorobenzene-d4	7.39

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1-Heptene	98	C7H14	000592-76-7	43
2		Octanal	128	C8H16O	000124-13-0	38
3		10-Azido-1-decanethiol	215	C10H21N3S	1152113-44-4	38
4		1-Hexene, 4-methyl-	98	C7H14	003769-23-1	35
5		(Z)-Hex-2-ene, 5-methyl-	98	C7H14	013151-17-2	30



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Data File : BN009831.D
Acq On : 22 Feb 2020 05:51
Operator : CG/JU
Sample : L1535-14
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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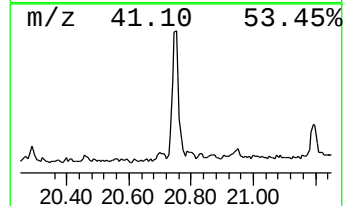
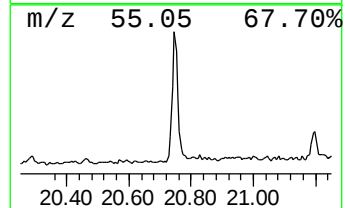
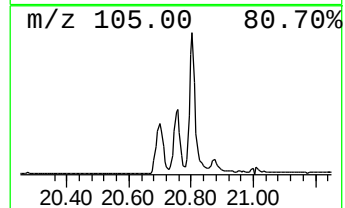
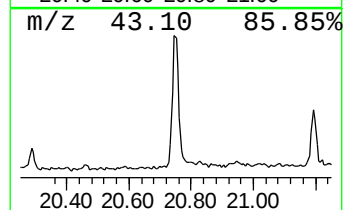
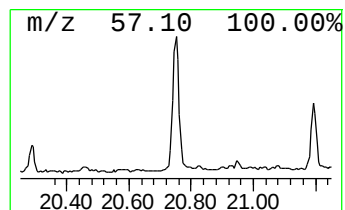
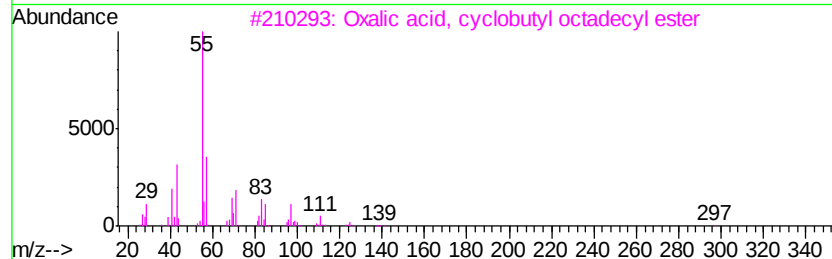
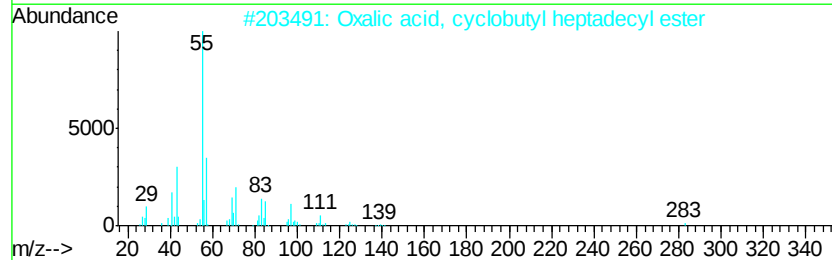
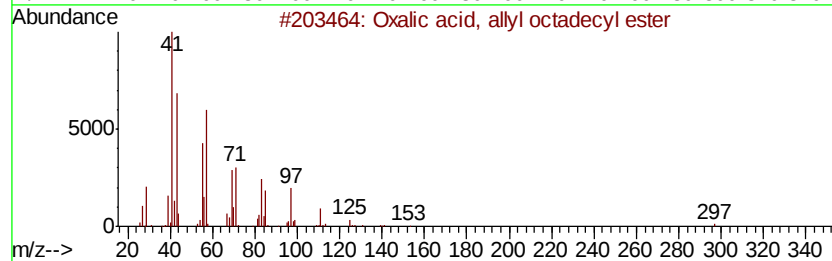
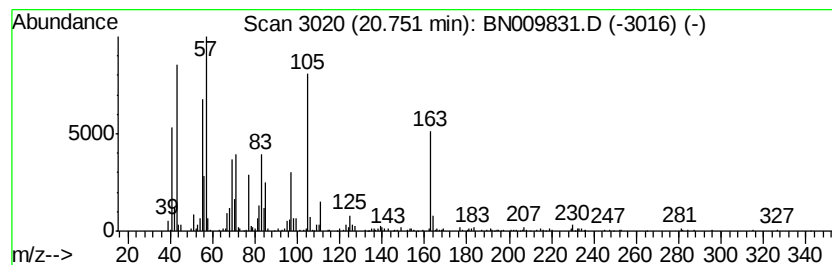
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 Oxalic acid, allyl octadecyl... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.75	6.50 ng/ul	768689	Chrysene-d12	21.00

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Oxalic acid, allyl octadecyl ester	382	C23H42O4	1000309-24-5	64
2		Oxalic acid, cyclobutyl heptadecyl ester	382	C23H42O4	1000309-70-7	58
3		Oxalic acid, cyclobutyl octadecyl ester	396	C24H44O4	1000309-70-8	58
4		Oxalic acid, isobutyl octadecyl ester	398	C24H46O4	1000309-38-3	58
5		Oxalic acid, allyl hexadecyl ester	354	C21H38O4	1000309-24-4	58



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN022220\
Data File : BN009831.D
Acq On : 22 Feb 2020 05:51
Operator : CG/JU
Sample : L1535-14
Misc :
ALS Vial : 20 Sample Multiplier: 1

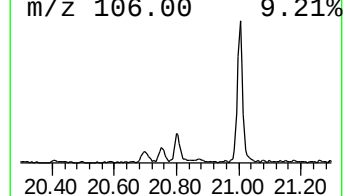
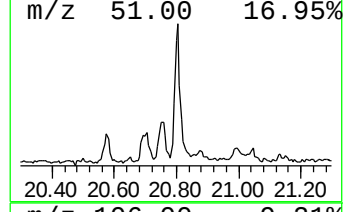
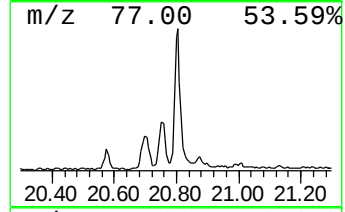
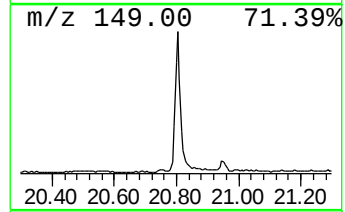
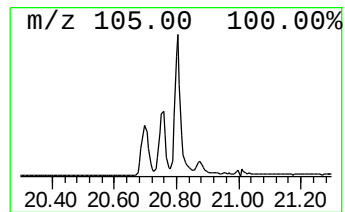
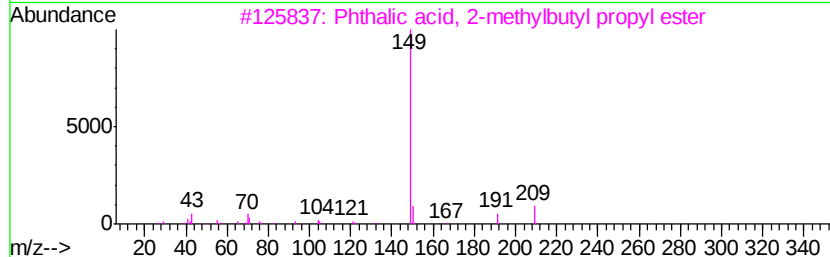
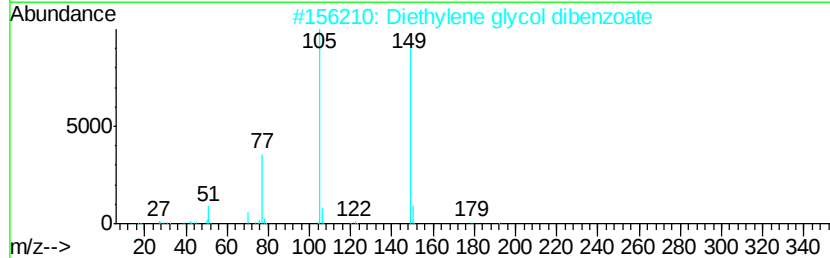
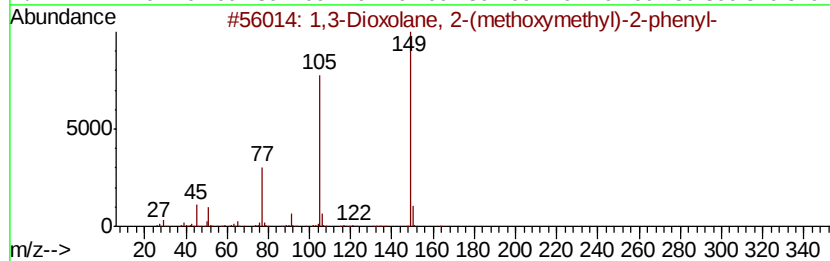
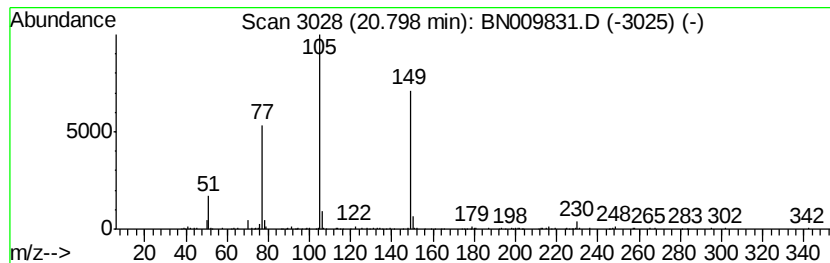
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 1,3-Dioxolane, 2-(methoxyme... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.	
20.80	3.76 ng/ul	444469	Chrysene-d12	21.00	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,3-Dioxolane, 2-(methoxymethyl)...	194	C11H14O3	1000156-71-2	64
2	Diethylene glycol dibenzoate	314	C18H18O5	000120-55-8	64
3	Phthalic acid, 2-methylbutyl pro...	278	C16H22O4	1000322-81-1	38
4	Phthalic acid, hexyl 2-methylbut...	320	C19H28O4	1000322-81-4	38
5	1,2-Bis(benzoyloxy)benzene	318	C20H14O4	000643-94-7	35



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN022220\
Data File : BN009831.D
Acq On : 22 Feb 2020 05:51
Operator : CG/JU
Sample : L1535-14
Misc :
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Instrument :
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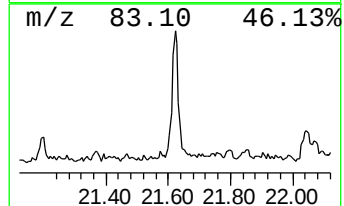
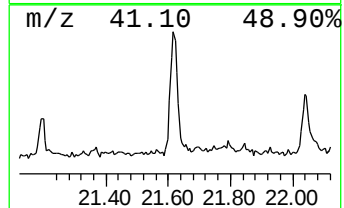
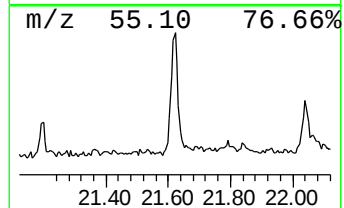
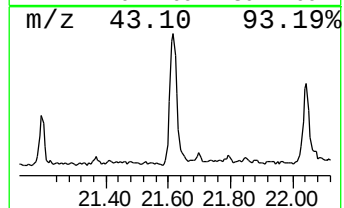
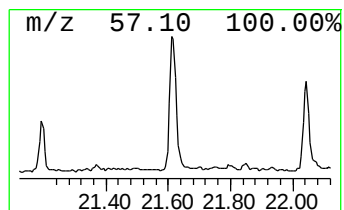
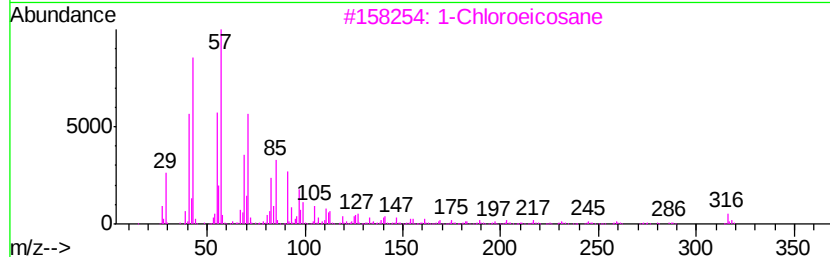
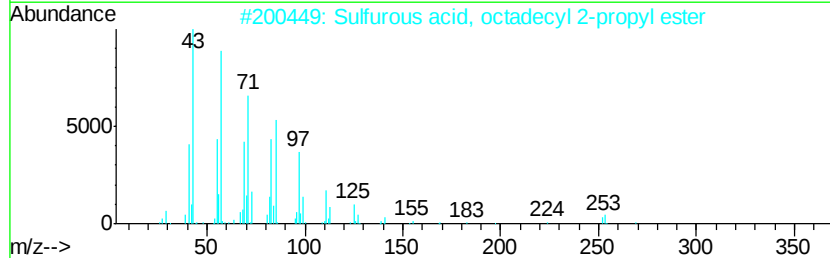
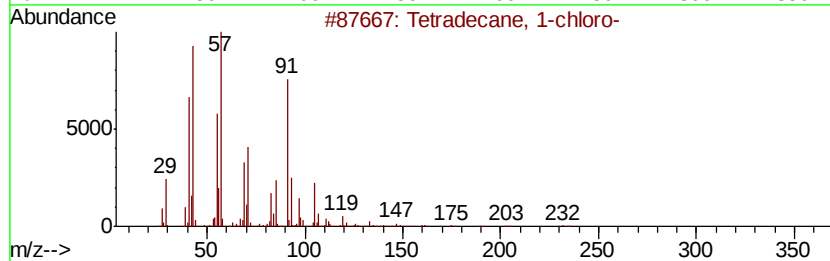
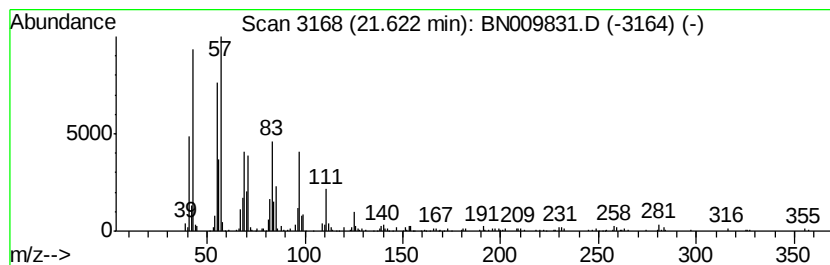
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 Tetradecane, 1-chloro- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.62	5.14 ng/ul	608851	Chrysene-d12	21.00

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tetradecane, 1-chloro-	232	C14H29Cl	002425-54-9	90
2		Sulfurous acid, octadecyl 2-propyl ester	376	C21H44O3S	1000309-12-7	81
3		1-Chloroeicosane	316	C20H41Cl	042217-02-7	72
4		Oxalic acid, pentadecyl propyl ester	342	C20H38O4	1000309-26-8	72
5		Nonadecane, 1-chloro-	302	C19H39Cl	062016-76-6	70



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN022220\
Data File : BN009831.D
Acq On : 22 Feb 2020 05:51
Operator : CG/JU
Sample : L1535-14
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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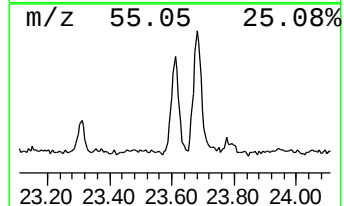
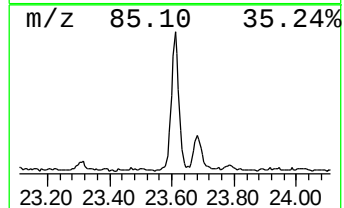
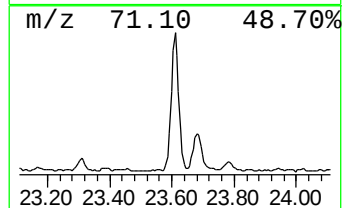
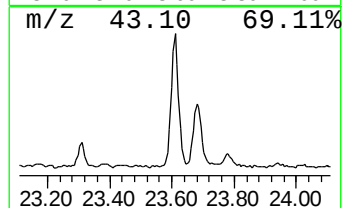
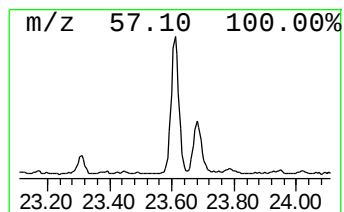
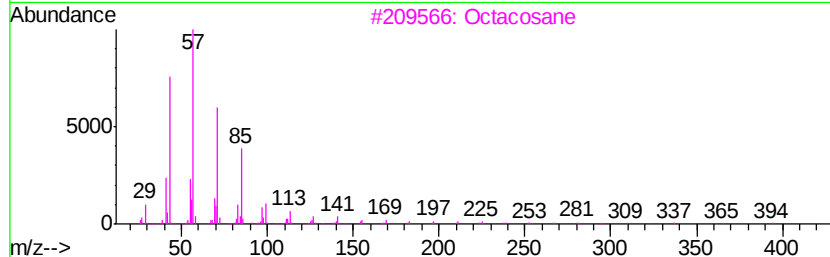
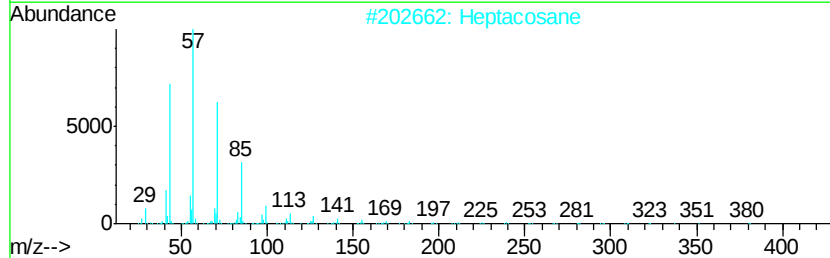
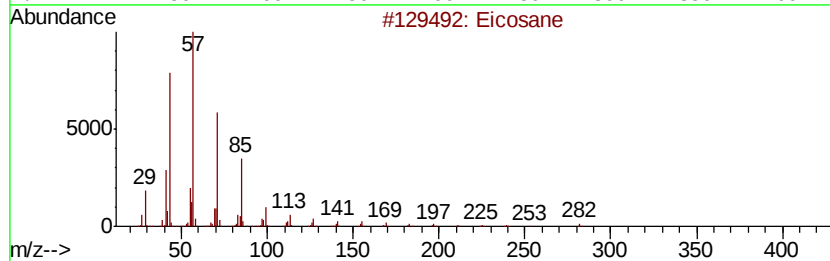
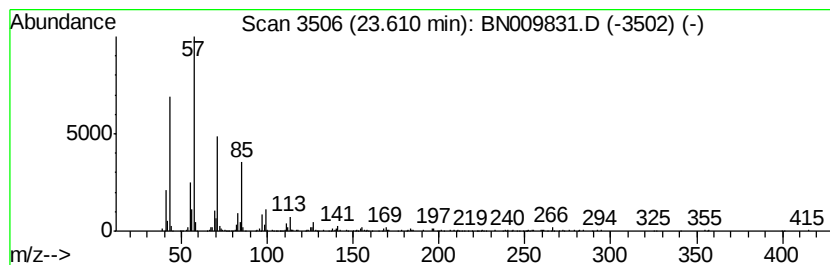
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 (DEL) Alkane: Straight-Chai... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.61	11.63 ng/ul	864327	Perylene-d12	23.10

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Eicosane	282	C20H42	000112-95-8	95
2		Heptacosane	380	C27H56	000593-49-7	90
3		Octacosane	394	C28H58	000630-02-4	90
4		Hexacosane	366	C26H54	000630-01-3	87
5		Hentriacontane	437	C31H64	000630-04-6	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN022220\
Data File : BN009831.D
Acq On : 22 Feb 2020 05:51
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Sample : L1535-14
Misc :
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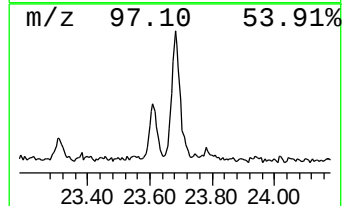
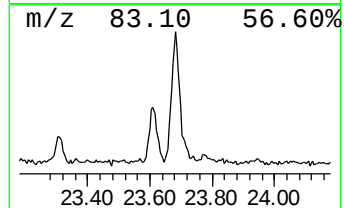
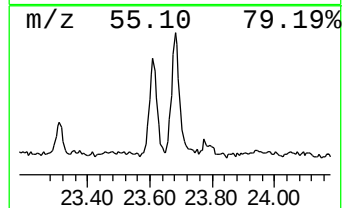
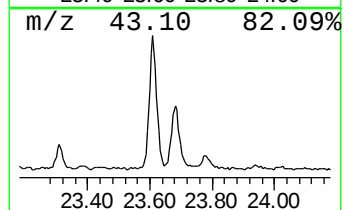
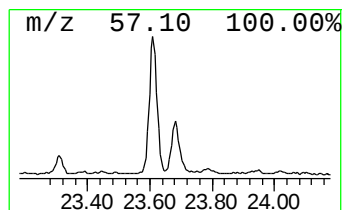
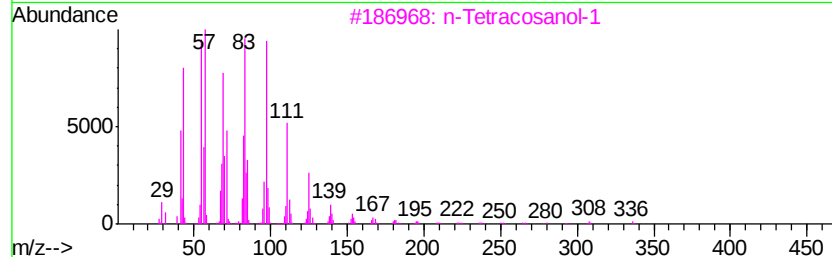
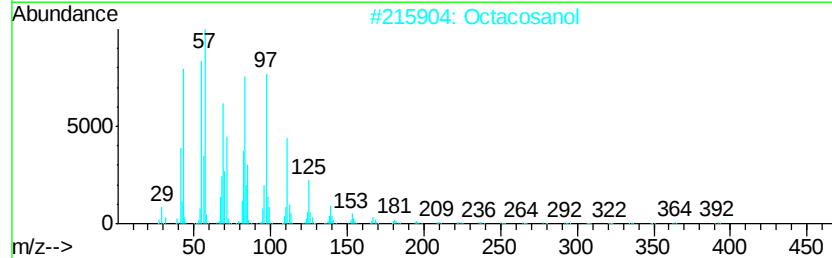
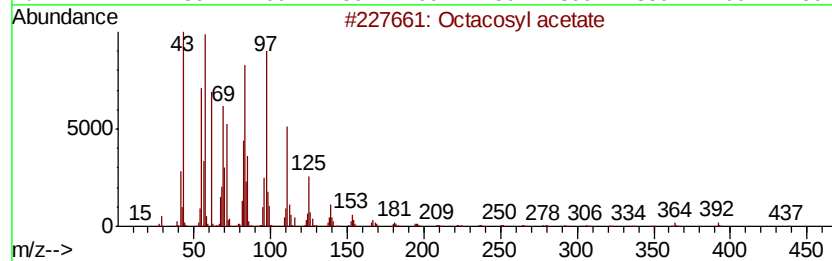
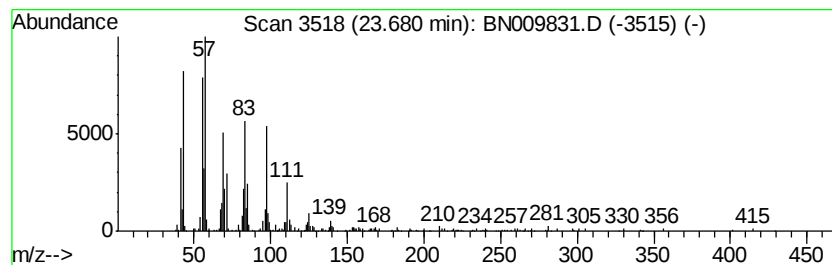
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 Octacosyl acetate Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.68	8.58 ng/ul	637549	Perylene-d12	23.10

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octacosyl acetate	452	C30H60O2	018206-97-8	91
2		Octacosanol	410	C28H58O	000557-61-9	91
3		n-Tetracosanol-1	354	C24H50O	000506-51-4	83
4		Triacontyl acetate	480	C32H64O2	041755-58-2	83
5		1-Pentadecene	210	C15H30	013360-61-7	68



Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN022220\
Data File : BN009831.D
Acq On : 22 Feb 2020 05:51
Operator : CG/JU
Sample : L1535-14
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument :
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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
Butane, 2,3-dimet...	3.93	3.2	ng/ul	121801	1	7.39	757998	20.0
unknown-01	4.33	2.9	ng/ul	108938	1	7.39	757998	20.0
(DEL) Alkane: Str...	8.71	2.4	ng/ul	89092	1	7.39	757998	20.0
Oxalic acid, ally...	20.75	6.5	ng/ul	768689	5	21.00	2366810	20.0
1,3-Dioxolane, 2-...	20.80	3.8	ng/ul	444469	5	21.00	2366810	20.0
Tetradecane, 1-ch...	21.62	5.1	ng/ul	608851	5	21.00	2366810	20.0
(DEL) Alkane: Str...	23.61	11.6	ng/ul	864327	6	23.10	1486260	20.0
Octacosyl acetate	23.68	8.6	ng/ul	637549	6	23.10	1486260	20.0