

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072421\
 Data File : BP006327.D
 Acq On : 25 Jul 2021 03:37
 Operator : CG/JU
 Sample : M3084-02
 Misc :
 ALS Vial : 29 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0022

Integration Parameters: LSCINT.P

Integrator: RTE
 Smoothing : OFF Filtering: 5
 Sampling : 1 Min Area: 0.1 % of largest Peak
 Start Thrs: 0.2 Max Peaks: 100
 Stop Thrs : 0 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
 Peak separation: 5

Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP072421.M
 Title : SVOA CALIBRATION

Signal : TIC: BP006327.D

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.311	33	38	48	rVB	66623	103104	1.27%	0.121%
2	3.711	102	106	122	rBV	366511	659190	8.09%	0.771%
3	3.934	140	144	149	rVB2	15230	22708	0.28%	0.027%
4	4.028	155	160	170	rVB	33424	55977	0.69%	0.065%
5	5.893	473	477	489	rVB	98562	157105	1.93%	0.184%
6	6.458	568	573	577	rBV	24111	34323	0.42%	0.040%
7	6.787	622	629	633	rVV6	9536	18533	0.23%	0.022%
8	6.964	651	659	670	rVB	336973	570216	7.00%	0.667%
9	7.134	681	688	705	rVB	1111764	1763621	21.65%	2.064%
10	7.322	712	720	731	rBV	1186411	1815818	22.29%	2.125%
11	7.422	731	737	742	rBV8	10789	18346	0.23%	0.021%
12	7.787	792	799	806	rBV	947647	1525123	18.72%	1.785%
13	7.858	806	811	817	rVB	95679	147393	1.81%	0.172%
14	8.034	834	841	846	rBV	46318	78245	0.96%	0.092%
15	8.411	900	905	909	rVV	11220	17599	0.22%	0.021%
16	8.493	911	919	929	rVV	1024386	1581429	19.41%	1.850%
17	8.581	929	934	936	rVV	37033	54389	0.67%	0.064%
18	8.611	936	939	946	rVB	40702	65664	0.81%	0.077%
19	8.881	979	985	990	rBV	316943	513033	6.30%	0.600%
20	8.946	990	996	1002	rVB	1274257	2087797	25.63%	2.443%
21	9.040	1007	1012	1021	rVV3	72939	151786	1.86%	0.178%
22	9.116	1021	1025	1030	rVB2	56961	81874	1.01%	0.096%
23	9.510	1085	1092	1097	rBV3	15931	31683	0.39%	0.037%
24	9.658	1110	1117	1130	rBV	985745	1669506	20.49%	1.953%
25	9.840	1136	1148	1154	rBV2	246290	621991	7.64%	0.728%
26	9.893	1154	1157	1165	rVB5	18543	34910	0.43%	0.041%
27	10.081	1178	1189	1197	rBV	327441	550899	6.76%	0.645%
28	10.193	1201	1208	1220	rVB	1581741	2589894	31.79%	3.030%
29	10.363	1231	1237	1246	rBV2	155275	243503	2.99%	0.285%
30	10.440	1246	1250	1254	rVV6	18735	30110	0.37%	0.035%
31	10.487	1254	1258	1265	rVB3	44457	80900	0.99%	0.095%
32	10.575	1265	1273	1279	rBV	1279140	2178255	26.74%	2.549%
33	10.634	1279	1283	1289	rVB3	61927	112776	1.38%	0.132%
34	10.710	1289	1296	1309	rBV	1564234	2641060	32.42%	3.090%
35	10.875	1320	1324	1331	rBV9	9784	18895	0.23%	0.022%
36	10.975	1337	1341	1347	rVB9	9237	19309	0.24%	0.023%

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Integration Parameters: LSCINT.P

Integrator: RTE

Soothing : OFF

Filtering: 5

Sampling : 1

Min Area: 0.1 % of largest Peak

Start Thrs: 0.2

Max Peaks: 100

Stop Thrs : 0

Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >

Peak separation: 5

Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP072421.M
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37	11.122	1357	1366	1371	rBV	181671	279095	3.43%	0.327%
38	11.181	1371	1376	1381	rVV	195940	311896	3.83%	0.365%
39	11.222	1381	1383	1389	rVB	36040	52945	0.65%	0.062%
40	11.369	1404	1408	1413	rBV5	14510	26137	0.32%	0.031%
41	11.752	1466	1473	1476	rBV	14918	24389	0.30%	0.029%
42	12.004	1507	1516	1522	rVB5	25859	54727	0.67%	0.064%
43	12.163	1534	1543	1548	rBV2	31392	57198	0.70%	0.067%
44	12.781	1643	1648	1655	rBV	74407	112311	1.38%	0.131%
45	13.034	1686	1691	1697	rBV	113727	159365	1.96%	0.186%
46	13.263	1725	1730	1735	rVB	33590	47097	0.58%	0.055%
47	13.340	1735	1743	1747	rBV	122369	192519	2.36%	0.225%
48	13.663	1791	1798	1802	rBV2	47138	76903	0.94%	0.090%
49	13.834	1815	1827	1843	rBV	3308324	4624108	56.77%	5.411%
50	13.957	1843	1848	1852	rVB5	25580	41039	0.50%	0.048%
51	14.104	1862	1873	1882	rBV	3493728	5144755	63.16%	6.020%
52	14.281	1898	1903	1906	rVB	15063	19812	0.24%	0.023%
53	14.328	1907	1911	1917	rVB4	14019	20195	0.25%	0.024%
54	14.416	1917	1926	1934	rBV	1871997	2822108	34.64%	3.302%
55	14.504	1936	1941	1945	rVB2	28401	40826	0.50%	0.048%
56	14.610	1950	1959	1967	rBV	377156	545249	6.69%	0.638%
57	14.840	1993	1998	2003	rBV4	29640	43760	0.54%	0.051%
58	14.928	2010	2013	2016	rBV3	12169	17776	0.22%	0.021%
59	15.016	2024	2028	2032	rBV3	14978	19907	0.24%	0.023%
60	15.104	2037	2043	2048	rVV4	19333	35840	0.44%	0.042%
61	15.169	2051	2054	2058	rVV	12882	18683	0.23%	0.022%
62	15.228	2060	2064	2070	rVV2	28170	56221	0.69%	0.066%
63	15.281	2070	2073	2078	rVB3	12575	17160	0.21%	0.020%
64	15.410	2088	2095	2102	rBV	4812775	6595480	80.97%	7.717%
65	15.522	2107	2114	2130	rBV	1497171	2089056	25.65%	2.444%
66	15.651	2130	2136	2141	rBV	54974	82237	1.01%	0.096%
67	15.781	2152	2158	2164	rBV2	36350	50381	0.62%	0.059%
68	15.969	2187	2190	2198	rBV9	8487	21838	0.27%	0.026%
69	16.098	2207	2212	2218	rBV7	19682	44832	0.55%	0.052%
70	16.204	2226	2230	2237	rVB5	40297	64089	0.79%	0.075%
71	16.316	2243	2249	2254	rBV6	13966	33456	0.41%	0.039%
72	16.410	2261	2265	2271	rVB3	14897	21563	0.26%	0.025%
73	16.510	2275	2282	2287	rBV10	7934	18010	0.22%	0.021%
74	16.569	2287	2292	2298	rVB2	34644	55285	0.68%	0.065%
75	16.669	2305	2309	2314	rVB3	22685	33588	0.41%	0.039%
76	16.910	2345	2350	2354	rBV	44108	73065	0.90%	0.085%

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77	16.945	2354	2356	2360	rVB4	18665	21233	0.26%	0.025%
78	17.163	2386	2393	2403	rBV	2228952	3162042	38.82%	3.700%
79	17.263	2404	2410	2417	rBV2	5219599	7412770	91.00%	8.673%
80	17.851	2504	2510	2517	rVV	2281640	2795249	34.31%	3.271%
81	17.969	2526	2530	2535	rVB3	34414	48053	0.59%	0.056%
82	18.075	2542	2548	2556	rBV	563024	779264	9.57%	0.912%
83	18.157	2556	2562	2570	rVB2	41402	114263	1.40%	0.134%
84	18.269	2578	2581	2585	rBV2	37799	47335	0.58%	0.055%
85	18.310	2586	2588	2592	rVB3	22039	22748	0.28%	0.027%
86	18.645	2642	2645	2648	rBV2	24159	26947	0.33%	0.032%
87	18.804	2669	2672	2676	rVB	29554	30524	0.37%	0.036%
88	18.892	2683	2687	2690	rBV2	31498	43242	0.53%	0.051%
89	19.081	2715	2719	2724	rBV2	68908	91211	1.12%	0.107%
90	19.151	2727	2731	2736	rVV2	78658	149597	1.84%	0.175%
91	19.192	2736	2738	2744	rVB4	35686	51317	0.63%	0.060%
92	19.310	2753	2758	2767	rBV	387494	523015	6.42%	0.612%
93	19.445	2778	2781	2785	rVB	48015	53891	0.66%	0.063%
94	19.504	2785	2791	2798	rBV	6207137	8146046	100.00%	9.531%
95	20.833	3015	3017	3025	rVB2	65158	66626	0.82%	0.078%
96	20.986	3040	3043	3049	rBV	339501	404527	4.97%	0.473%
97	21.263	3085	3090	3098	rBV	2521120	3277416	40.23%	3.835%
98	22.369	3274	3278	3283	rBV3	259779	384612	4.72%	0.450%
99	23.457	3456	3463	3472	rVB2	4203422	8088781	99.30%	9.464%
100	23.604	3482	3488	3499	rVB2	1656496	3332181	40.91%	3.899%

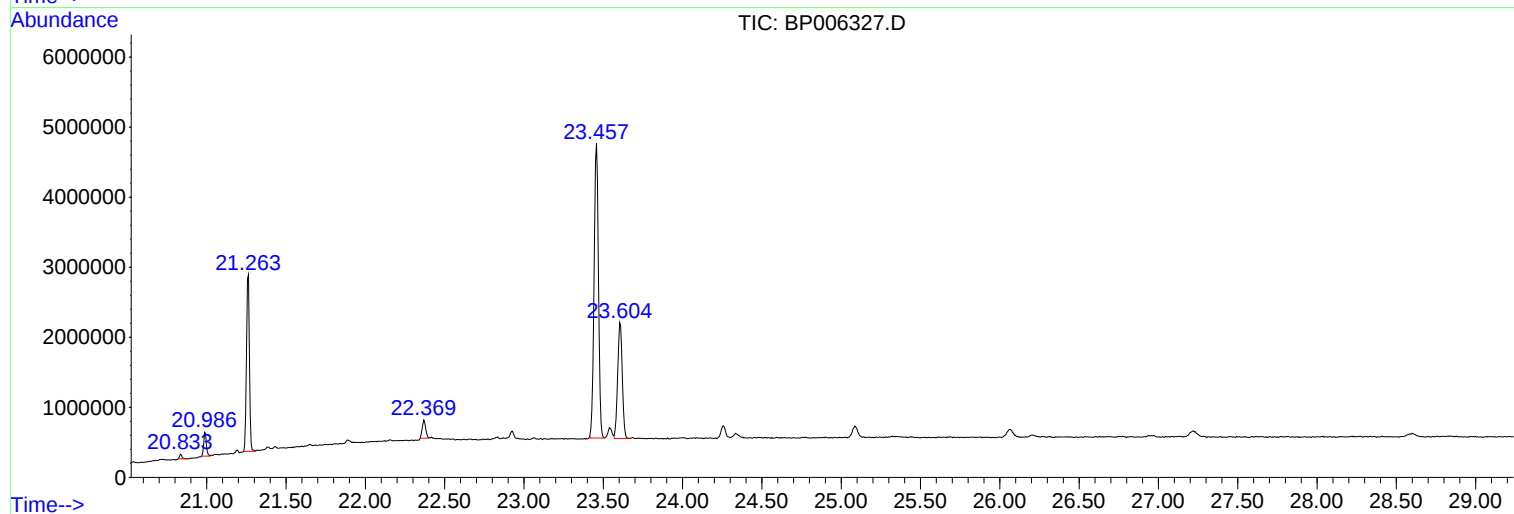
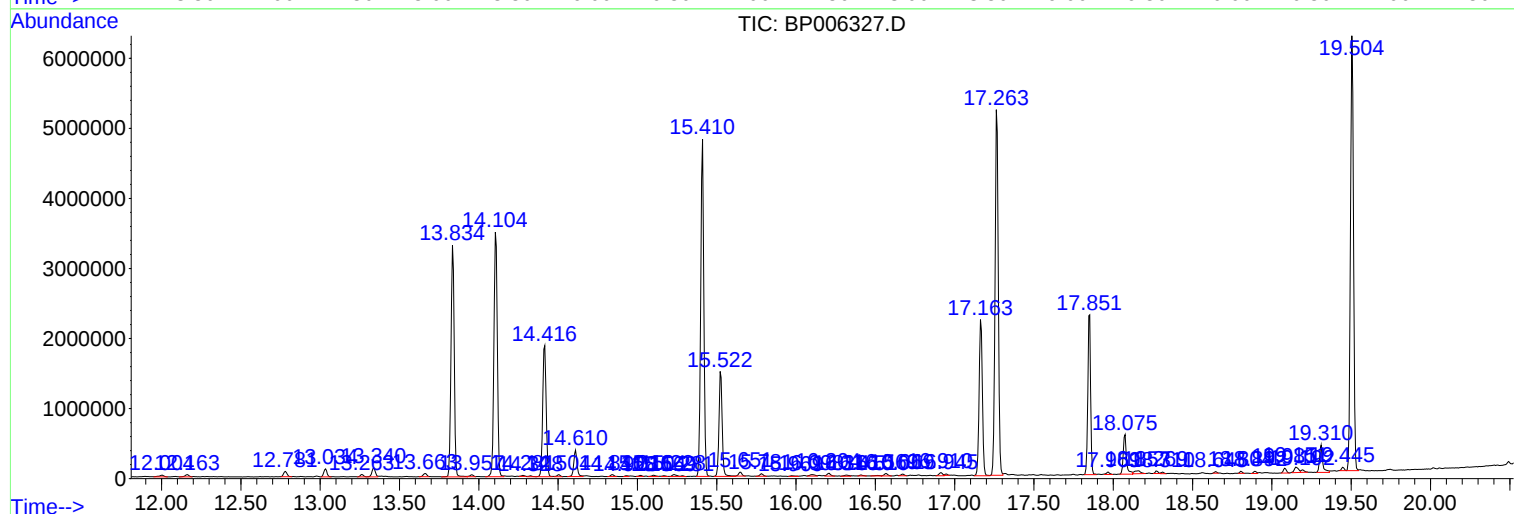
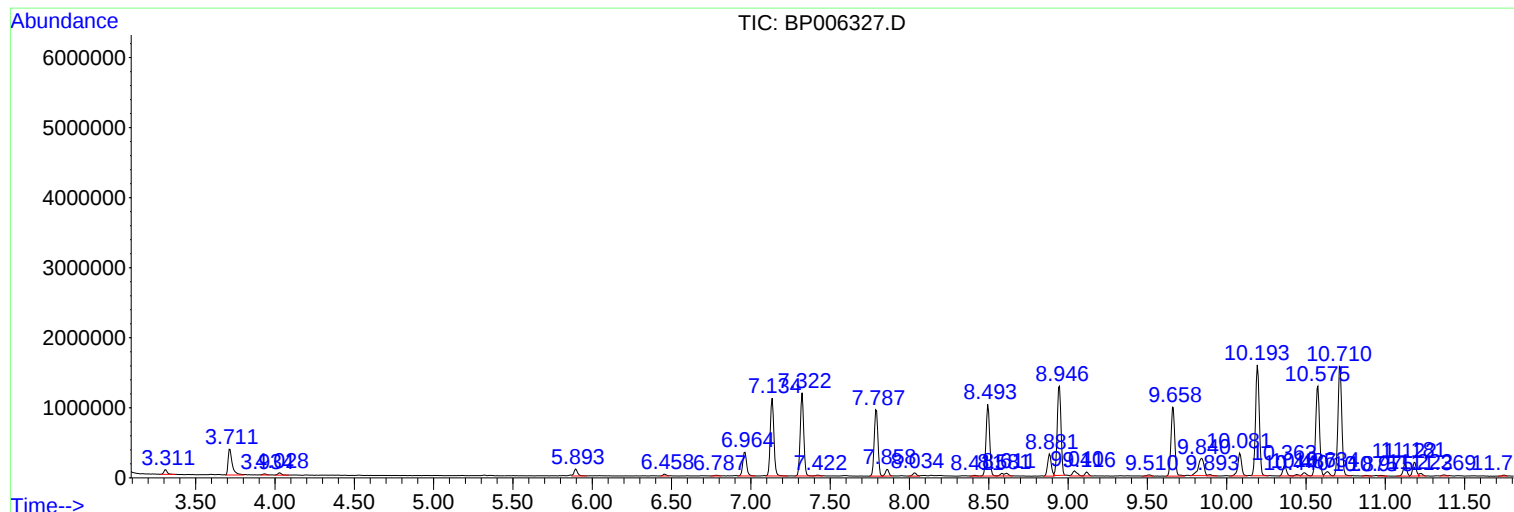
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Misc :
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Instrument :
BNA_P
ClientSampleId :
C0022

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP072421.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P



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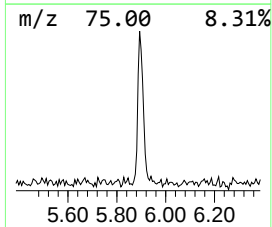
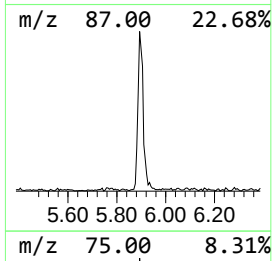
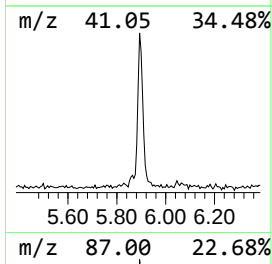
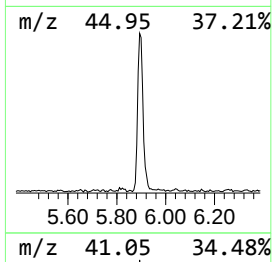
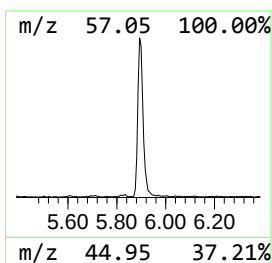
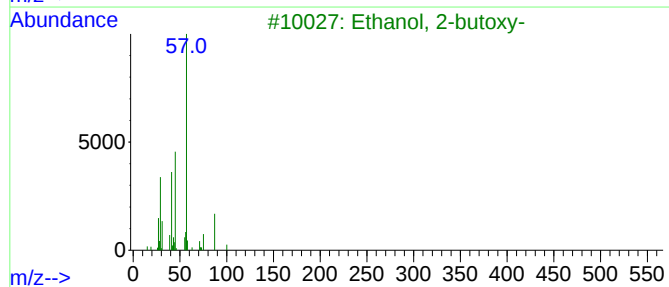
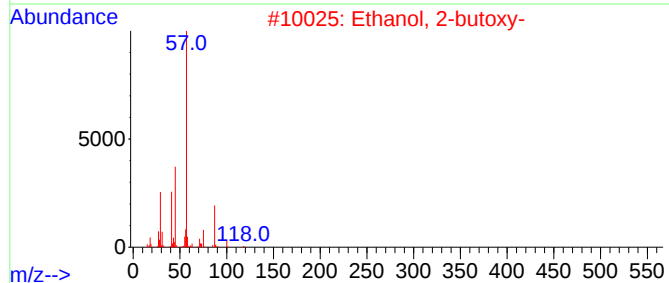
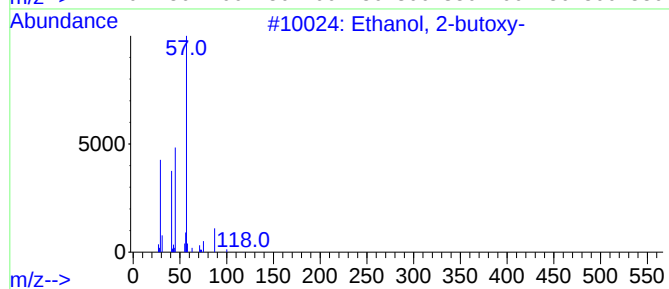
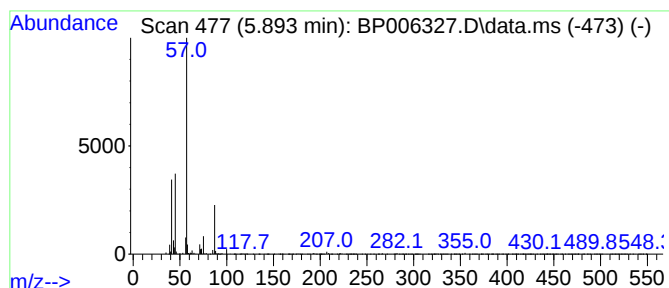
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP072421.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 Ethanol, 2-butoxy- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.890	2.06 ng/ul	157105	1,4-Dichlorobenzene-d4	7.793

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Ethanol, 2-butoxy-	118	C6H14O2	000111-76-2	59
2			Ethanol, 2-butoxy-	118	C6H14O2	000111-76-2	59
3			Ethanol, 2-butoxy-	118	C6H14O2	000111-76-2	50
4			Ethanol, 2-butoxy-	118	C6H14O2	000111-76-2	50
5			Ethylene glycol monoisobutyl ether	118	C6H14O2	004439-24-1	47



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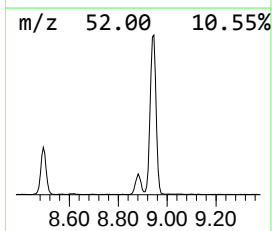
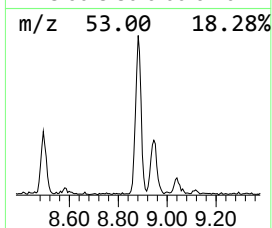
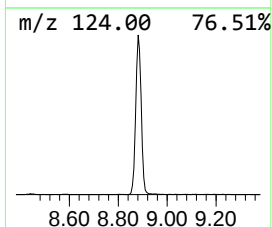
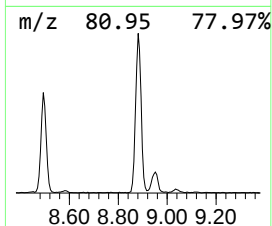
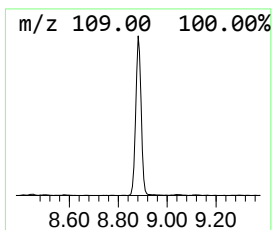
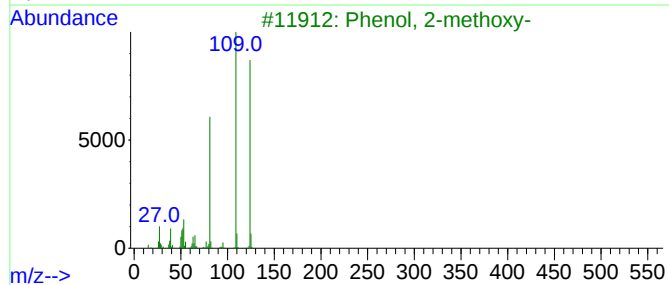
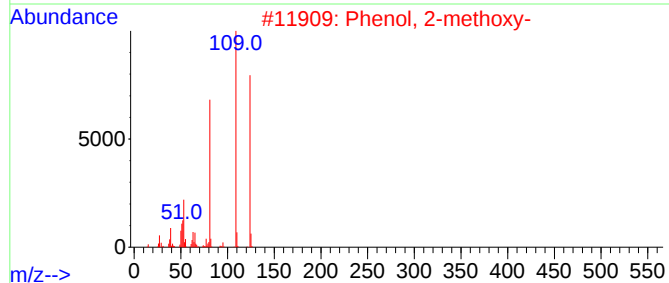
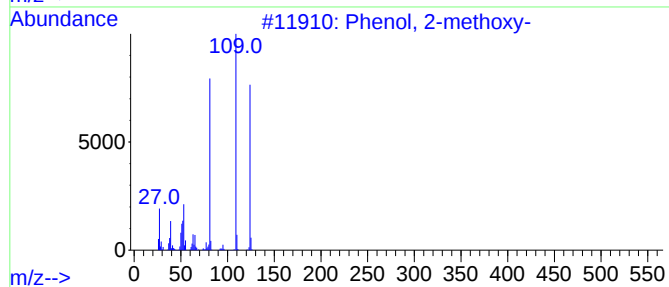
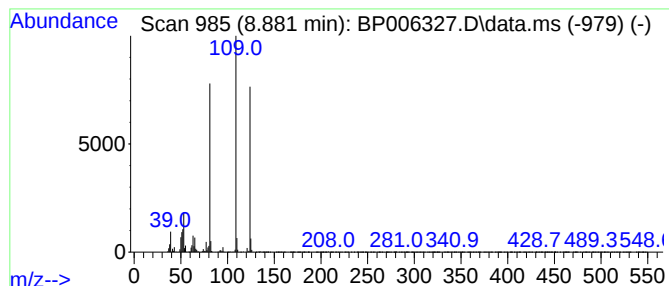
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 Phenol, 2-methoxy- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.880	6.73 ng/ul	513033	1,4-Dichlorobenzene-d4	7.793

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, 2-methoxy-		124	C7H8O2	000090-05-1	97	
2	Phenol, 2-methoxy-		124	C7H8O2	000090-05-1	96	
3	Phenol, 2-methoxy-		124	C7H8O2	000090-05-1	70	
4	Phenol, 2-methoxy-		124	C7H8O2	000090-05-1	70	
5	Mequinol		124	C7H8O2	000150-76-5	64	



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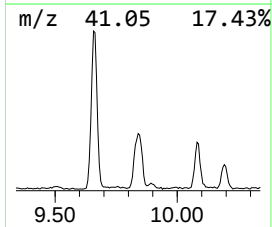
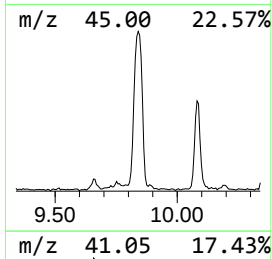
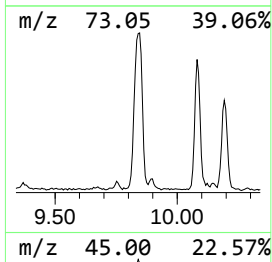
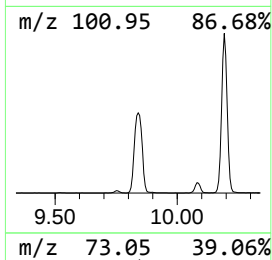
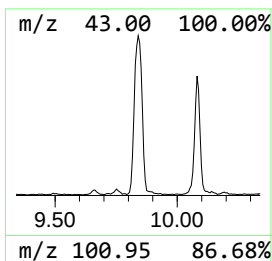
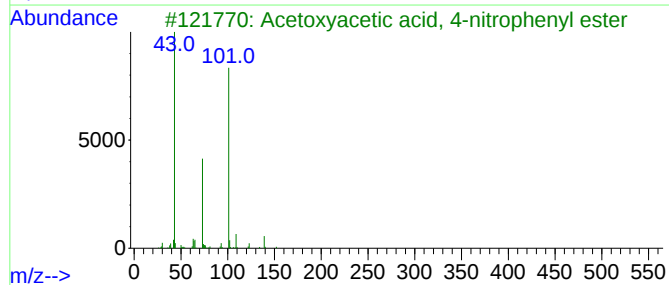
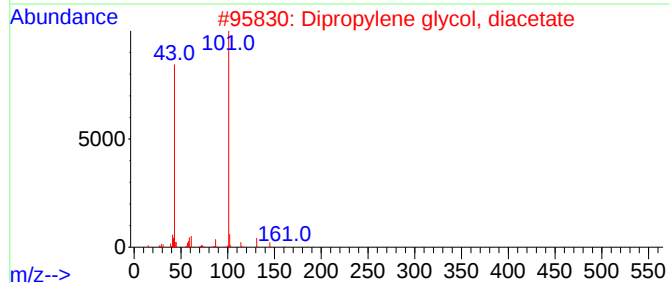
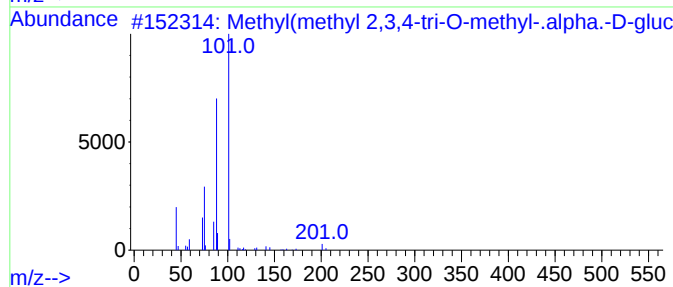
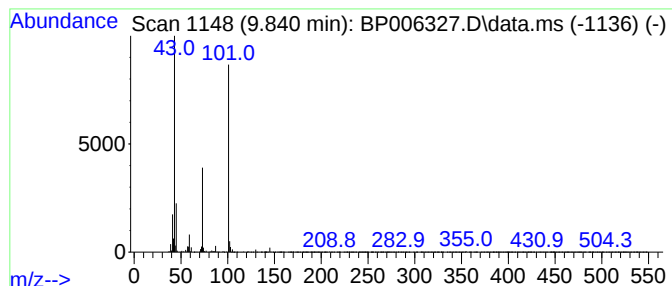
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TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 Methyl(methyl 2,3,4-tri-O-m... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.840	5.71 ng/ul	621991	Naphthalene-d8	10.575

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Methyl(methyl 2,3,4-tri-O-methyl...	264	C11H20O7	052729-97-2	64
2			Dipropylene glycol, diacetate	218	C10H18O5	1000506-25-8	59
3			Acetoxyacetic acid, 4-nitropheny...	239	C10H9NO6	1000307-54-5	45
4			Adipic acid, hexyl 4-methoxy-2-m...	330	C18H34O5	1000324-29-9	42
5			3H-1,2,4-Triazole-3-thione, 1,2-...	101	C2H3N3S	003179-31-5	42



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072421\
Data File : BP006327.D
Acq On : 25 Jul 2021 03:37
Operator : CG/JU
Sample : M3084-02
Misc :
ALS Vial : 29 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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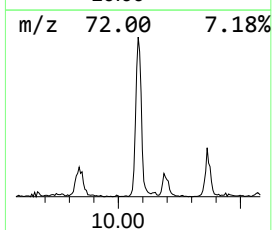
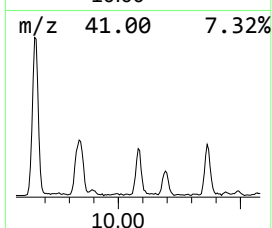
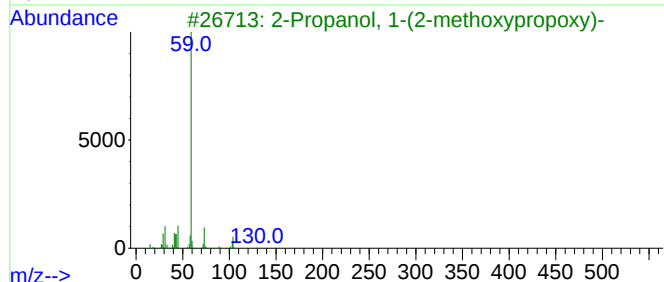
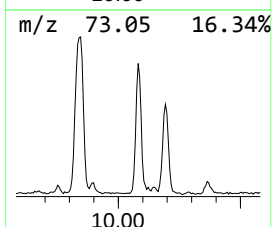
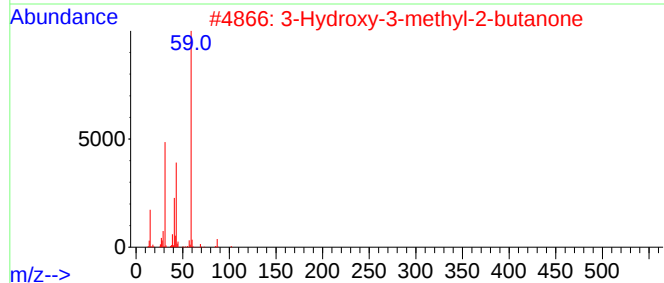
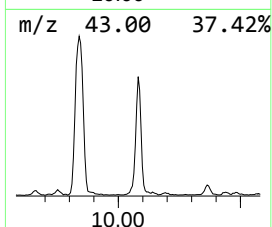
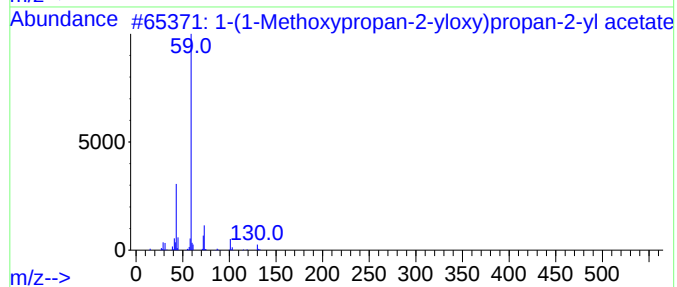
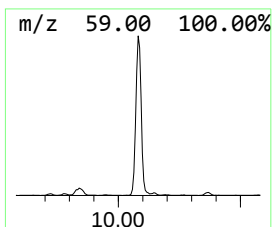
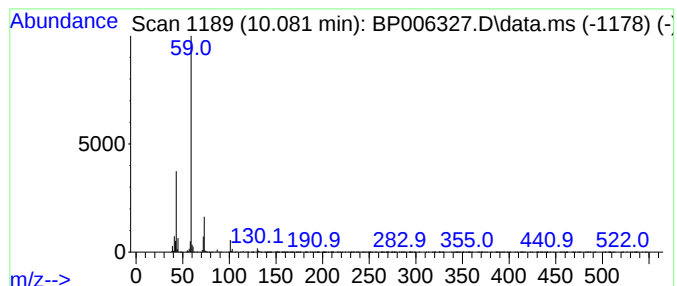
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 1-(1-Methoxypropan-2-yloxy)... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.080	5.06 ng/ul	550899	Naphthalene-d8	10.575

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-(1-Methoxypropan-2-yloxy)propa...	190	C9H18O4	1000367-08-6	90		
2	3-Hydroxy-3-methyl-2-butanone	102	C5H10O2	000115-22-0	53		
3	2-Propanol, 1-(2-methoxypropoxy)-	148	C7H16O3	013429-07-7	50		
4	3-Hexanol	102	C6H14O	000623-37-0	45		
5	2-Hexanol, methyl ether	116	C7H16O	1000333-92-0	45		



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072421\
Data File : BP006327.D
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Sample : M3084-02
Misc :
ALS Vial : 29 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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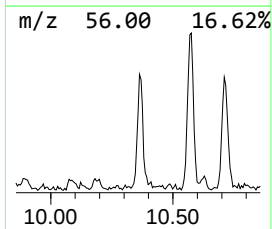
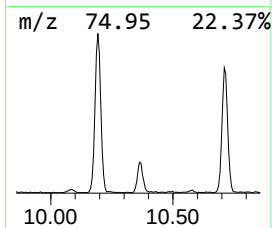
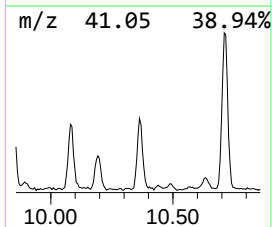
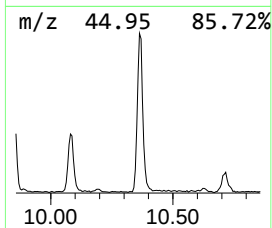
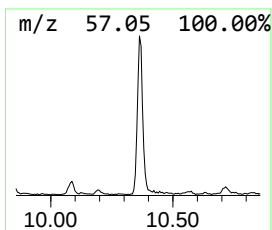
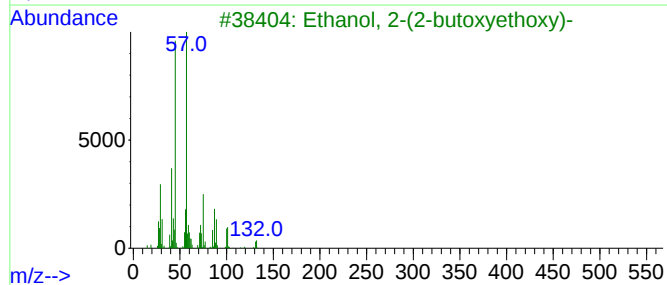
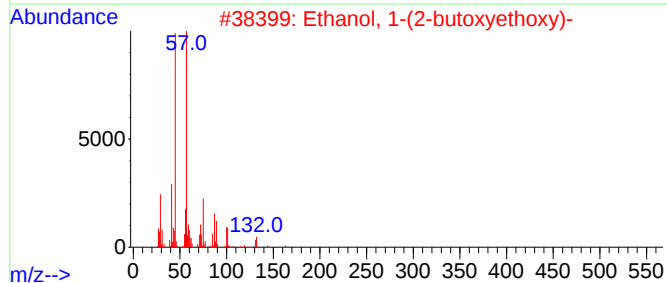
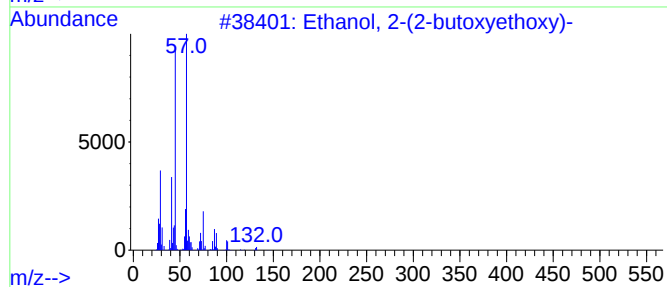
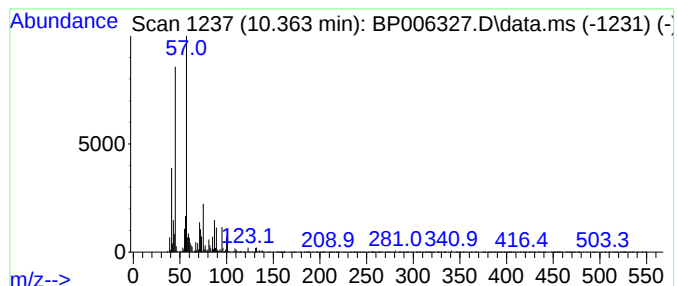
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 Ethanol, 2-(2-butoxyethoxy)- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.360	2.24 ng/ul	243503	Naphthalene-d8	10.575

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	90
2			Ethanol, 1-(2-butoxyethoxy)-	162	C8H18O3	054446-78-5	90
3			Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	83
4			Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	83
5			Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072421\
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Sample : M3084-02
Misc :
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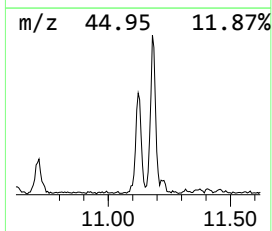
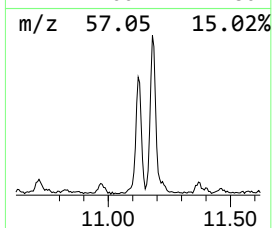
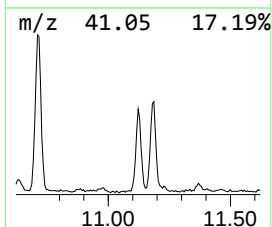
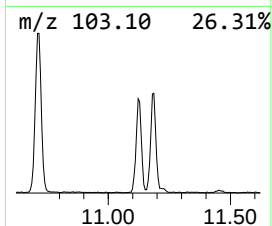
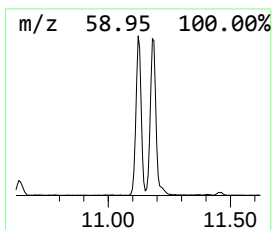
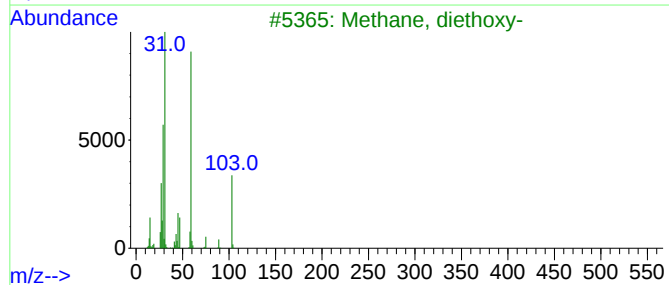
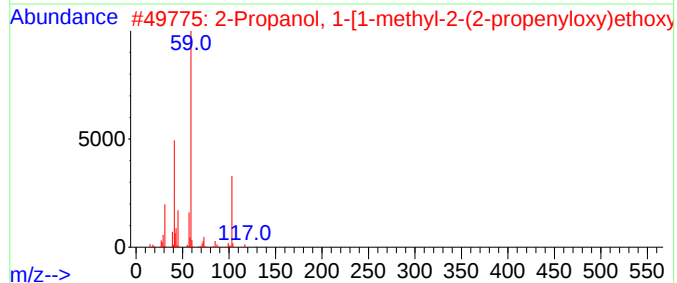
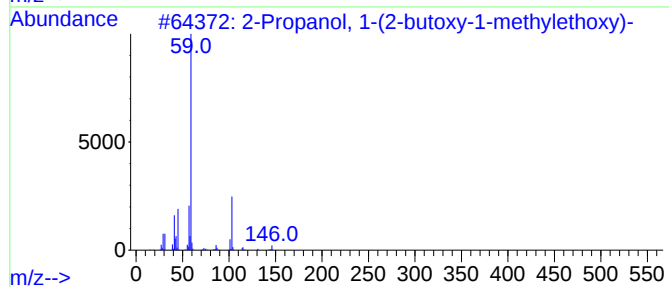
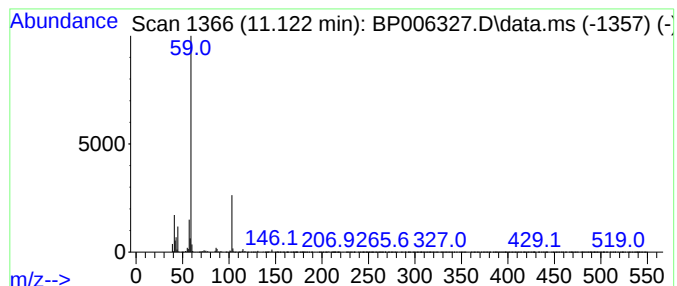
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 2-Propanol, 1-(2-butoxy-1-m... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.120	2.56 ng/ul	279095	Naphthalene-d8	10.575

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Propanol, 1-(2-butoxy-1-methyl...	190	C10H22O3	029911-28-2	83		
2	2-Propanol, 1-[1-methyl-2-(2-pro...	174	C9H18O3	055956-25-7	83		
3	Methane, diethoxy-	104	C5H12O2	000462-95-3	80		
4	2-Propanol, 1-(2-butoxy-1-methyl...	190	C10H22O3	029911-28-2	74		
5	Methane, diethoxy-	104	C5H12O2	000462-95-3	72		



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072421\
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Sample : M3084-02
Misc :
ALS Vial : 29 Sample Multiplier: 1

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ClientSampleId :
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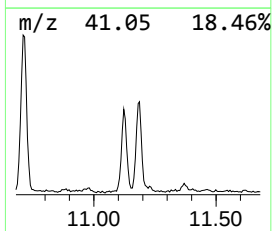
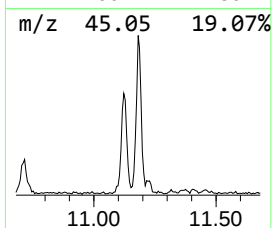
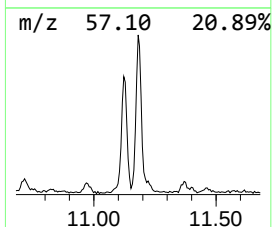
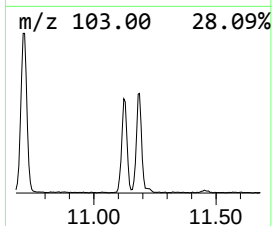
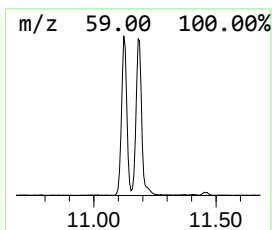
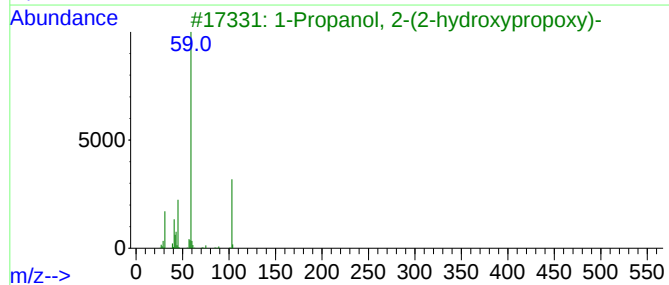
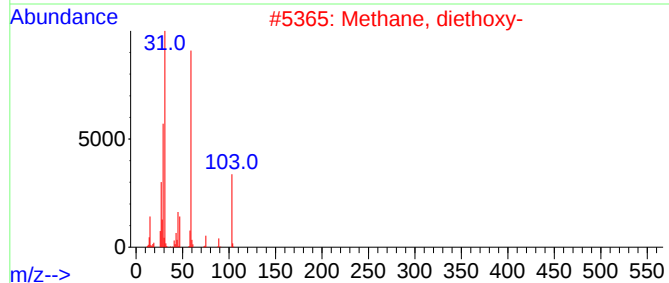
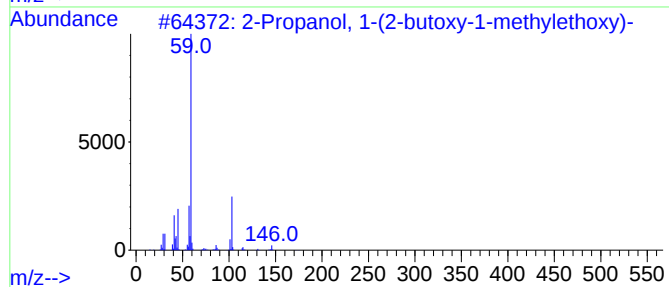
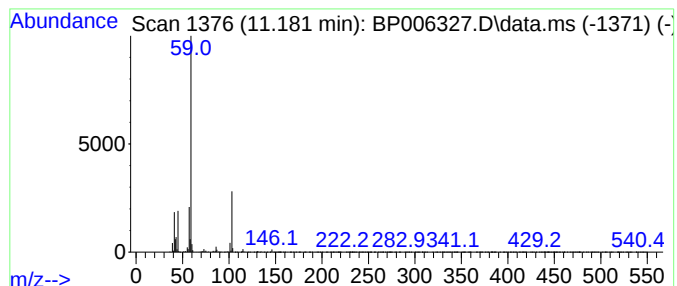
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 7 Methane, diethoxy- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.180	2.86 ng/ul	311896	Naphthalene-d8	10.575

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Propanol, 1-(2-butoxy-1-methyl...	190	C10H22O3	029911-28-2	90		
2	Methane, diethoxy-	104	C5H12O2	000462-95-3	78		
3	1-Propanol, 2-(2-hydroxypropoxy)-	134	C6H14O3	000106-62-7	78		
4	2-Propanol, 1,1'-[(1-methyl-1,2-...	192	C9H20O4	001638-16-0	64		
5	Dipropylene glycol (isomer 2)	134	C6H14O3	1000506-25-4	64		



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072421\
Data File : BP006327.D
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Sample : M3084-02
Misc :
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Instrument :
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ClientSampleId :
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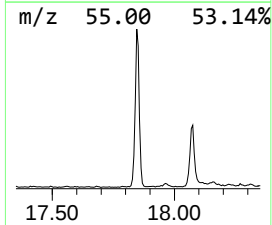
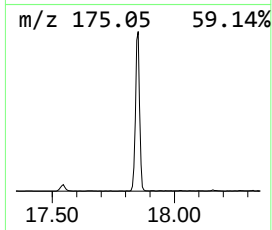
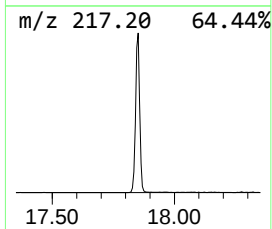
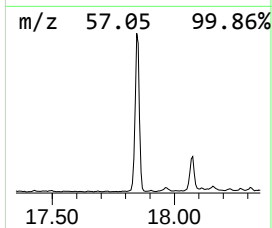
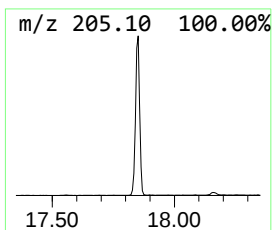
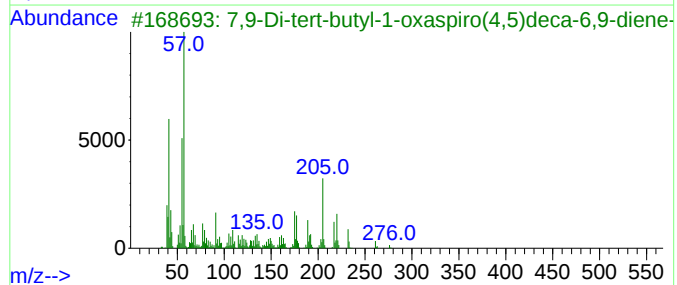
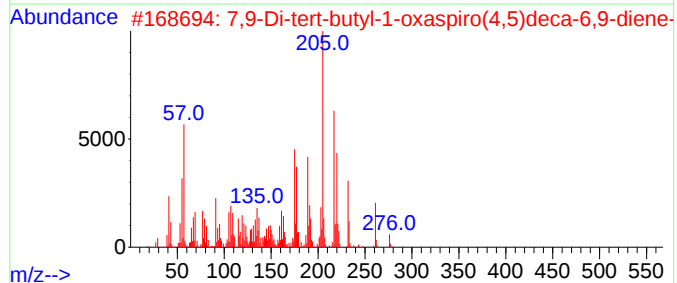
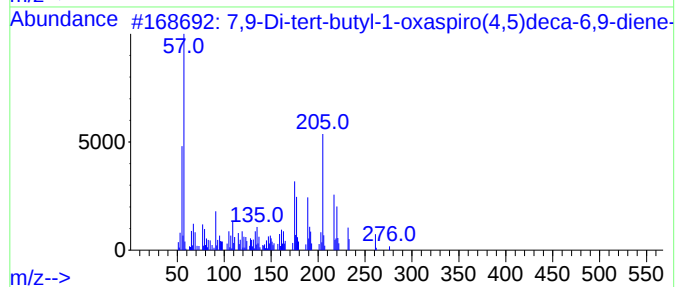
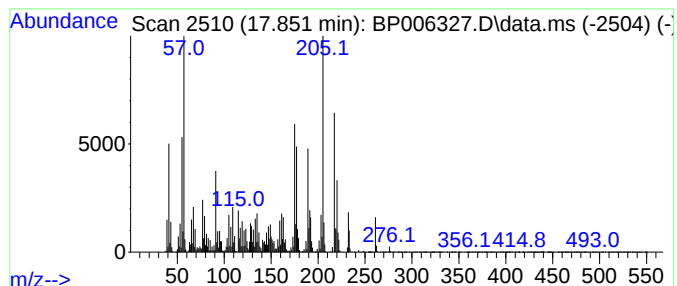
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 8 7,9-Di-tert-butyl-1-oxaspiro... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.850	17.68 ng/ul	2795250	Phenanthrene-d10	17.163

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	7,9-Di-tert-butyl-1-oxaspiro(4,5...	276	C17H24O3	082304-66-3	99		
2	7,9-Di-tert-butyl-1-oxaspiro(4,5...	276	C17H24O3	082304-66-3	99		
3	7,9-Di-tert-butyl-1-oxaspiro(4,5...	276	C17H24O3	082304-66-3	97		
4	2,5-Cyclohexadien-1-one, 2,6-bis...	232	C16H24O	006738-27-8	93		
5	Ethanone, 1-(5,6,7,8-tetrahydro-...	220	C14H20O2	071596-88-8	50		



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072421\
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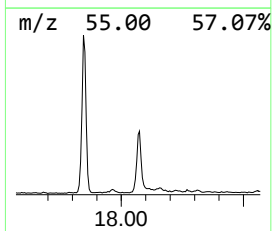
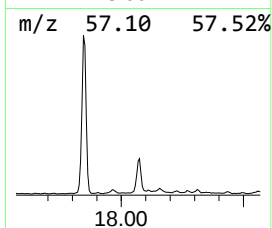
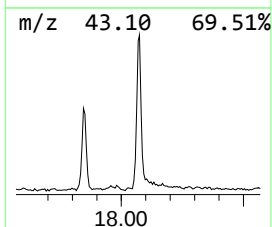
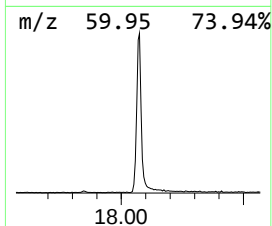
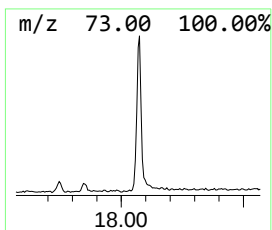
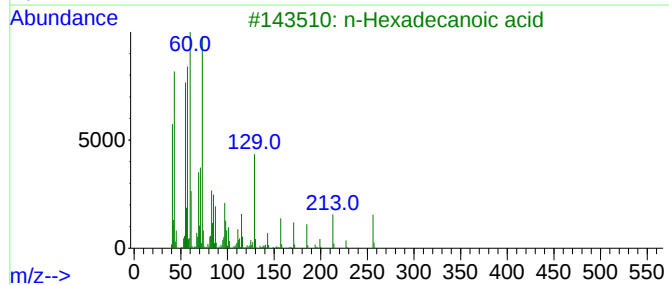
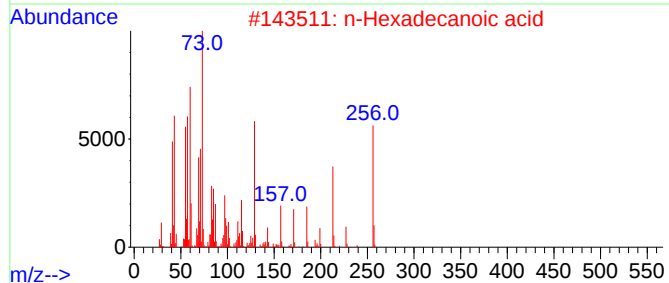
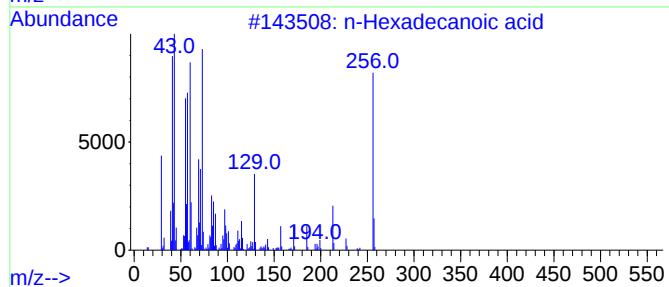
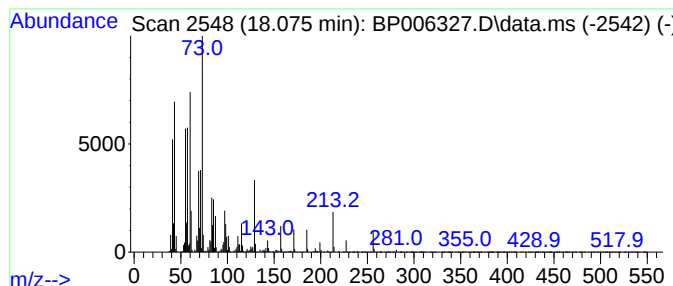
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 9 n-Hexadecanoic acid Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.070	4.93 ng/ul	779264	Phenanthrene-d10	17.163

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
3			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
4			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
5			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	96



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072421\
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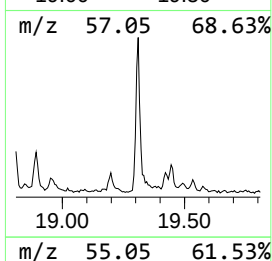
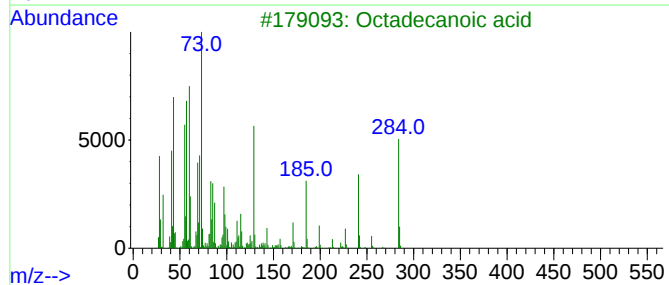
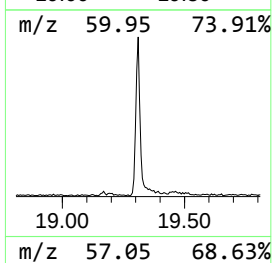
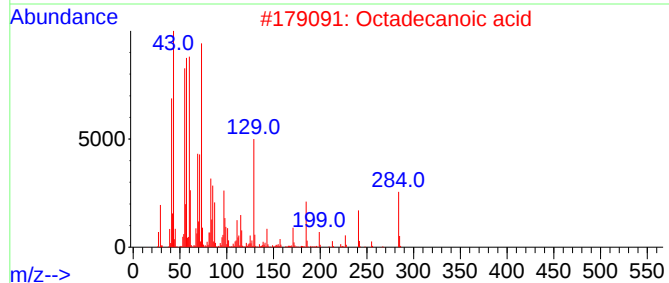
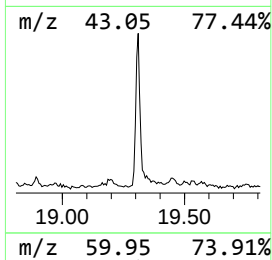
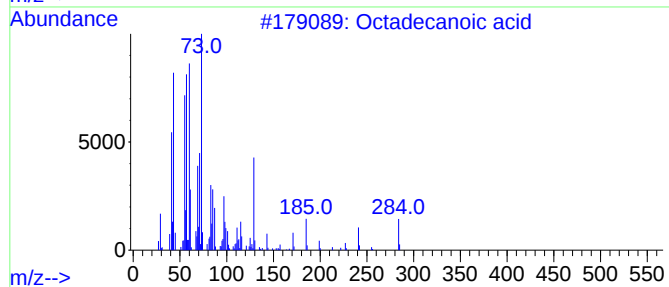
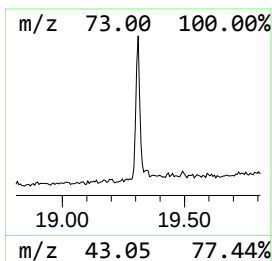
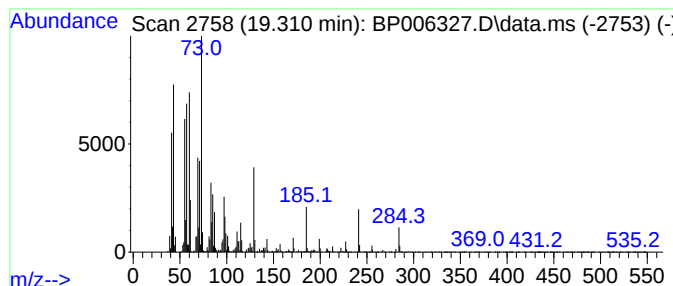
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 10 Octadecanoic acid Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.310	3.19 ng/ul	523015	Chrysene-d12	21.263

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octadecanoic acid	284	C18H36O2	000057-11-4	99
2			Octadecanoic acid	284	C18H36O2	000057-11-4	99
3			Octadecanoic acid	284	C18H36O2	000057-11-4	99
4			Octadecanoic acid	284	C18H36O2	000057-11-4	94
5			Octadecanoic acid	284	C18H36O2	000057-11-4	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072421\
Data File : BP006327.D
Acq On : 25 Jul 2021 03:37
Operator : CG/JU
Sample : M3084-02
Misc :
ALS Vial : 29 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0022

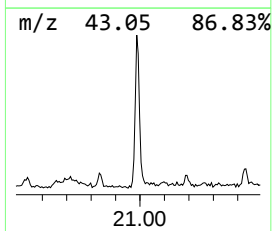
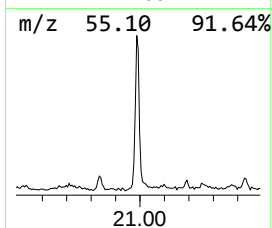
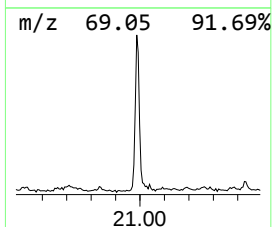
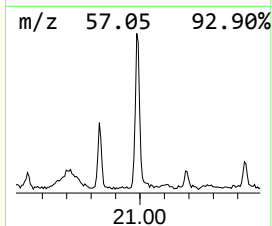
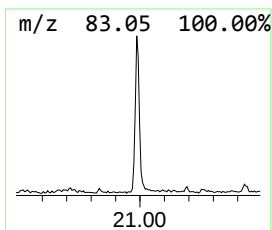
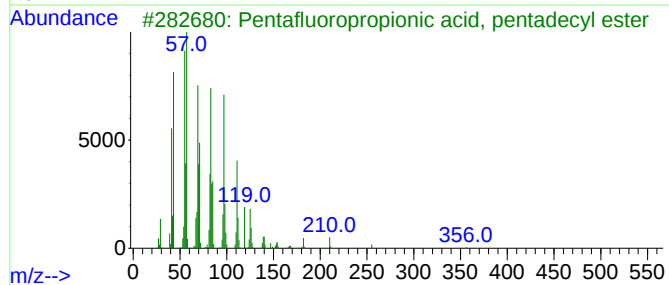
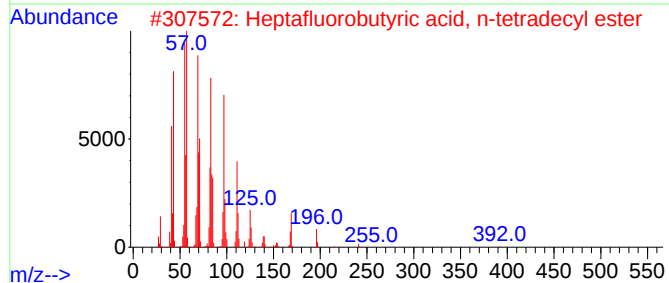
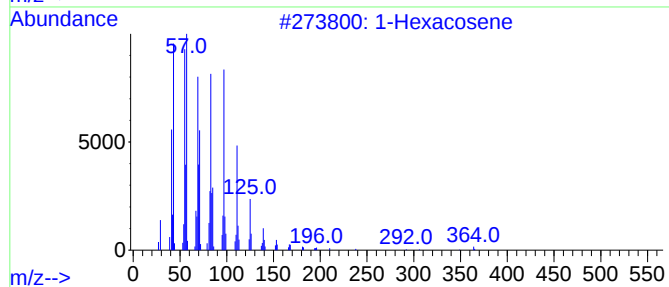
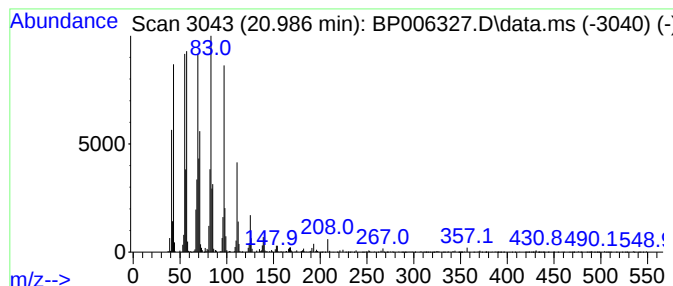
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 11 (DEL) Alkane: Straight-Chai... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.990	2.47 ng/ul	404527	Chrysene-d12	21.263

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Hexacosene			364	C26H52	018835-33-1	98
2	Heptafluorobutyric acid, n-tetra...			410	C18H29F7O2	007365-36-8	94
3	Pentafluoropropionic acid, penta...			374	C18H31F5O2	959092-08-1	94
4	Trifluoroacetic acid, pentadecyl...			324	C17H31F3O2	959010-23-2	94
5	Heptafluorobutyric acid, pentade...			424	C19H31F7O2	959261-23-5	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072421\
Data File : BP006327.D
Acq On : 25 Jul 2021 03:37
Operator : CG/JU
Sample : M3084-02
Misc :
ALS Vial : 29 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0022

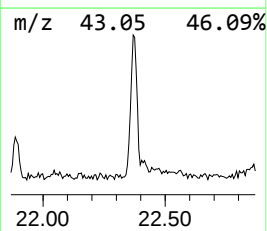
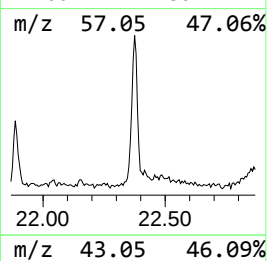
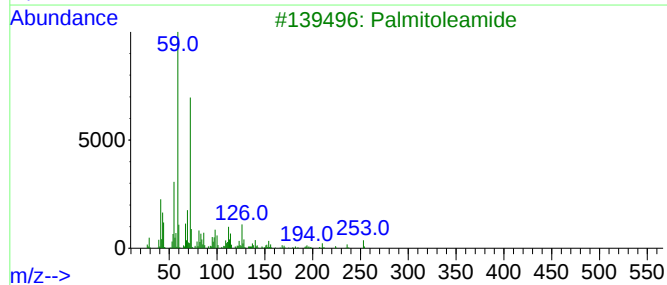
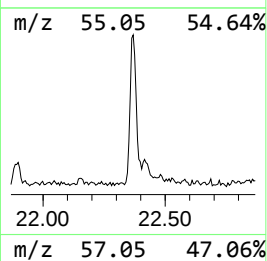
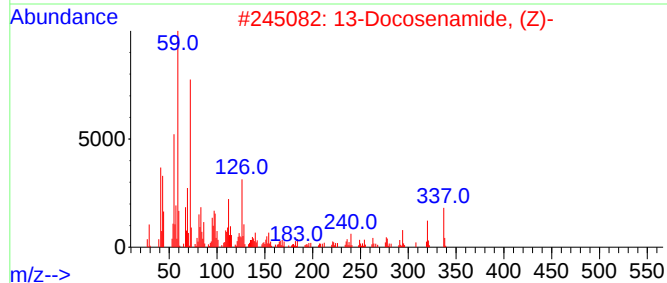
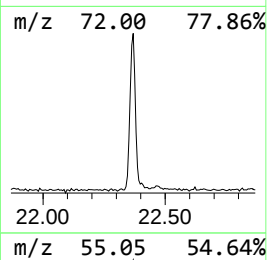
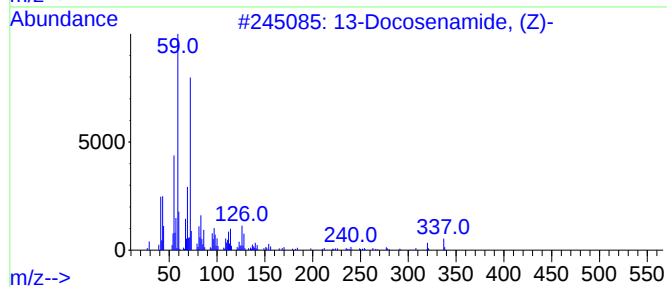
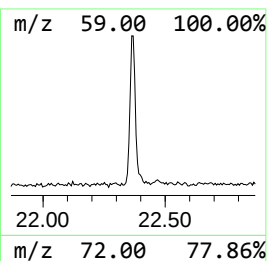
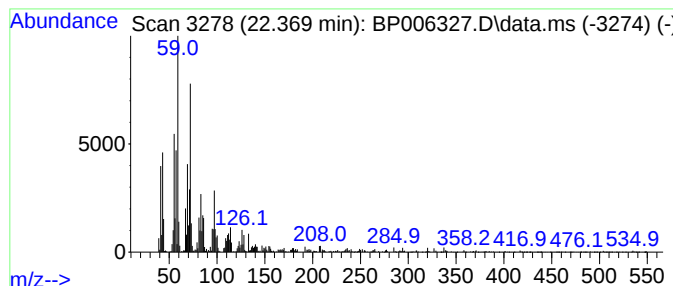
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 12 13-Docosenamide, (Z)- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.370	2.35 ng/ul	384612	Chrysene-d12	21.263

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		13-Docosenamide, (Z)-	337	C22H43NO	000112-84-5	98
2		13-Docosenamide, (Z)-	337	C22H43NO	000112-84-5	91
3		Palmitoleamide	253	C16H31NO	106010-22-4	83
4		9-Octadecenamide, (Z)-	281	C18H35NO	000301-02-0	80
5		9-Octadecenamide, (Z)-	281	C18H35NO	000301-02-0	58



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072421\
 Data File : BP006327.D
 Acq On : 25 Jul 2021 03:37
 Operator : CG/JU
 Sample : M3084-02
 Misc :
 ALS Vial : 29 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0022

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP072421.M
 Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST0.L
 TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Ethanol, 2-butoxy-	5.890	2.1	ng/ul	157105	1	7.793	1525120	20.0
Phenol, 2-methoxy-	8.880	6.7	ng/ul	513033	1	7.793	1525120	20.0
Methyl(methyl 2...	9.840	5.7	ng/ul	621991	2	10.575	2178260	20.0
1-(1-Methoxypro...	10.080	5.1	ng/ul	550899	2	10.575	2178260	20.0
Ethanol, 2-(2-b...	10.360	2.2	ng/ul	243503	2	10.575	2178260	20.0
2-Propanol, 1-(...	11.120	2.6	ng/ul	279095	2	10.575	2178260	20.0
Methane, diethoxy-	11.180	2.9	ng/ul	311896	2	10.575	2178260	20.0
7,9-Di-tert-but...	17.850	17.7	ng/ul	2795250	4	17.163	3162040	20.0
n-Hexadecanoic ...	18.070	4.9	ng/ul	779264	4	17.163	3162040	20.0
Octadecanoic acid	19.310	3.2	ng/ul	523015	5	21.263	3277420	20.0
(DEL) Alkane: S...	20.990	2.5	ng/ul	404527	5	21.263	3277420	20.0
13-Docosenamide...	22.370	2.4	ng/ul	384612	5	21.263	3277420	20.0