

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP080324\
 Data File : BP021340.D
 Acq On : 03 Aug 2024 11:38
 Operator : CG/JU
 Sample : P3278-02
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 E1ZP7

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-BP071024.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BP021340.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.164	27	30	36	rVB	24778	34740	1.04%	0.092%
2	3.464	78	81	93	rBV	38604	76846	2.30%	0.204%
3	3.587	100	102	126	rBV	219105	454966	13.63%	1.207%
4	3.881	149	152	160	rVB4	9041	15432	0.46%	0.041%
5	4.781	301	305	312	rVB2	5917	10378	0.31%	0.028%
6	6.858	650	658	674	rBV	250896	676211	20.26%	1.794%
7	6.981	674	679	696	rVB	292385	519334	15.56%	1.378%
8	7.181	707	713	730	rBV	349412	678799	20.33%	1.800%
9	7.316	731	736	745	rVB4	6928	21236	0.64%	0.056%
10	7.475	761	763	770	rVB8	2980	4752	0.14%	0.013%
11	7.628	782	789	801	rBV	585890	1024728	30.70%	2.718%
12	8.375	907	916	935	rBV	367041	879902	26.36%	2.334%
13	8.793	980	987	1004	rBV	343214	682606	20.45%	1.811%
14	9.358	1078	1083	1090	rBV4	9490	22826	0.68%	0.061%
15	9.505	1102	1108	1127	rBV	291406	598507	17.93%	1.588%
16	9.952	1179	1184	1192	rVB8	7237	16321	0.49%	0.043%
17	10.057	1195	1202	1221	rBV	271541	852827	25.55%	2.262%
18	10.381	1250	1257	1270	rBV	774488	1478715	44.30%	3.922%
19	10.563	1278	1288	1314	rBV	351252	911808	27.31%	2.419%
20	10.975	1354	1358	1369	rVB	6273	19195	0.58%	0.051%
21	11.175	1384	1392	1407	rBV4	48487	185231	5.55%	0.491%
22	11.593	1461	1463	1472	rVB10	2112	3978	0.12%	0.011%
23	12.004	1526	1533	1538	rBV8	6277	14556	0.44%	0.039%
24	12.581	1625	1631	1632	rBV6	2580	4242	0.13%	0.011%
25	12.846	1672	1676	1678	rBV5	2890	3906	0.12%	0.010%
26	13.675	1811	1817	1831	rBV	947440	1540006	46.13%	4.085%
27	13.951	1855	1864	1874	rBV	1149216	1798928	53.89%	4.772%
28	14.140	1894	1896	1902	rVB7	2650	3789	0.11%	0.010%
29	14.263	1907	1917	1929	rBV	1330367	2086594	62.51%	5.535%
30	14.475	1944	1953	1965	rBV5	51388	156418	4.69%	0.415%
31	14.628	1974	1979	1987	rBV	114548	341993	10.25%	0.907%
32	15.051	2048	2051	2056	rVV5	5222	7572	0.23%	0.020%
33	15.116	2060	2062	2069	rVB8	6070	9815	0.29%	0.026%
34	15.281	2082	2090	2111	rBV	1320656	2357799	70.63%	6.254%
35	15.475	2115	2123	2144	rVV	205079	602701	18.06%	1.599%
36	15.634	2147	2150	2152	rVB4	3096	3752	0.11%	0.010%

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37	15.845	2182	2186	2189	rBV6	6153	9370	0.28%	0.025%
38	16.004	2208	2213	2219	rBV7	6296	15469	0.46%	0.041%
39	16.175	2236	2242	2244	rBV6	5415	9402	0.28%	0.025%
40	16.822	2349	2352	2357	rBV6	7058	11329	0.34%	0.030%
41	16.987	2376	2380	2384	rBV7	4643	7421	0.22%	0.020%
42	17.081	2388	2396	2406	rBV	1538636	2623180	78.58%	6.958%
43	17.187	2407	2414	2428	rBV2	1614536	2585824	77.46%	6.859%
44	17.575	2478	2480	2484	rVB5	6443	7910	0.24%	0.021%
45	17.763	2508	2512	2519	rVB	28192	36788	1.10%	0.098%
46	17.869	2528	2530	2532	rBV3	4285	3943	0.12%	0.010%
47	18.010	2549	2554	2566	rBV	303789	546594	16.37%	1.450%
48	19.116	2739	2742	2748	rBV6	23460	36015	1.08%	0.096%
49	19.192	2752	2755	2759	rBV3	45592	77153	2.31%	0.205%
50	19.345	2777	2781	2787	rBV2	98996	168010	5.03%	0.446%
51	19.486	2803	2805	2808	rBV2	22590	24749	0.74%	0.066%
52	19.575	2814	2820	2826	rBV	2289837	3251970	97.42%	8.626%
53	20.128	2910	2914	2920	rVB	85089	108665	3.26%	0.288%
54	20.645	2999	3002	3007	rBV3	106417	154235	4.62%	0.409%
55	21.169	3086	3091	3102	rBV2	285659	587172	17.59%	1.557%
56	21.527	3146	3152	3165	rBV	1886300	2925497	87.64%	7.760%
57	24.586	3665	3672	3684	rVB	1181687	3070447	91.98%	8.144%
58	24.816	3704	3711	3726	rVB	1095398	3338136	100.00%	8.854%

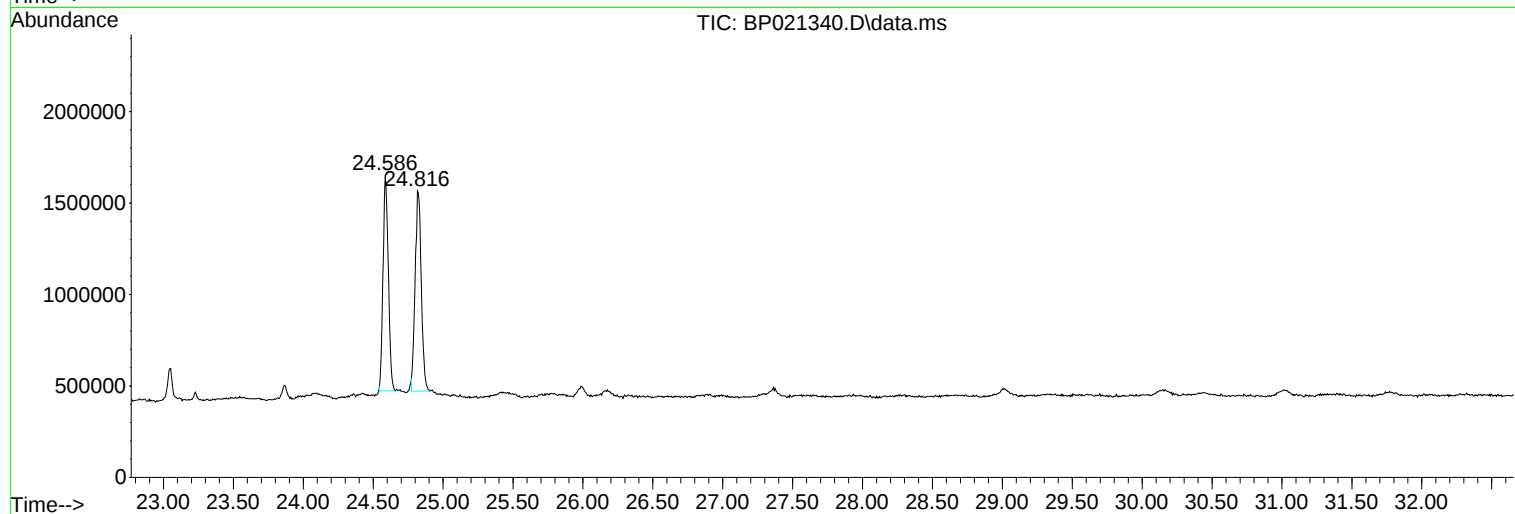
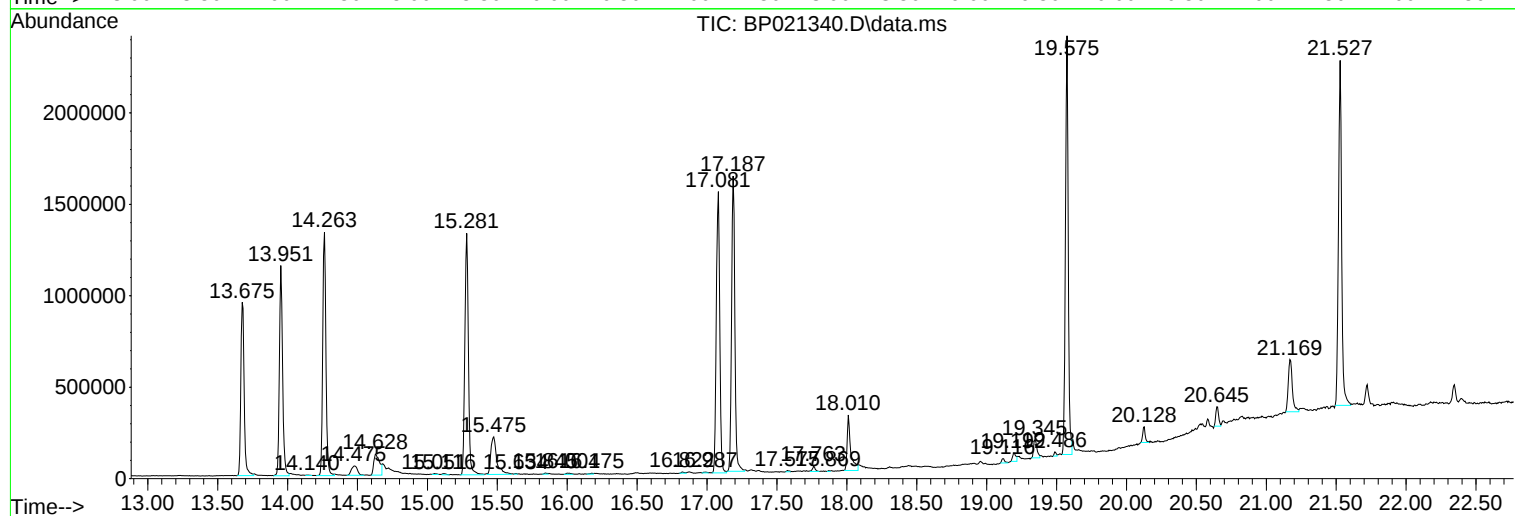
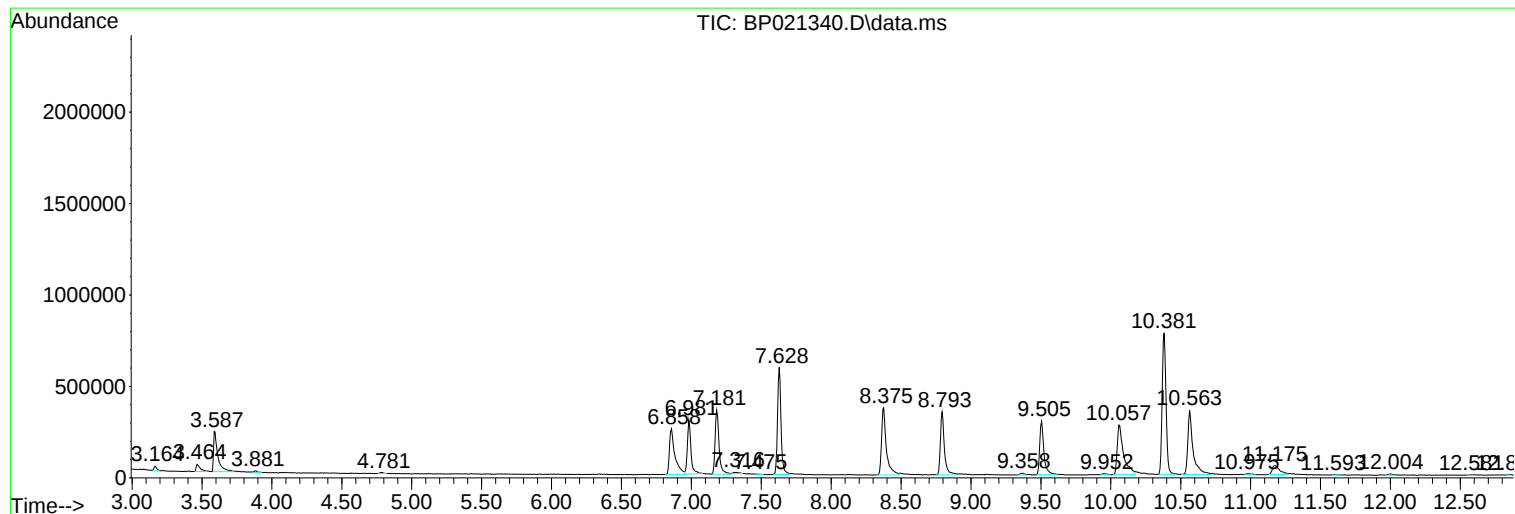
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Quant Title : SVOA CALIBRATION

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



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ALS Vial : 5 Sample Multiplier: 1

Instrument :
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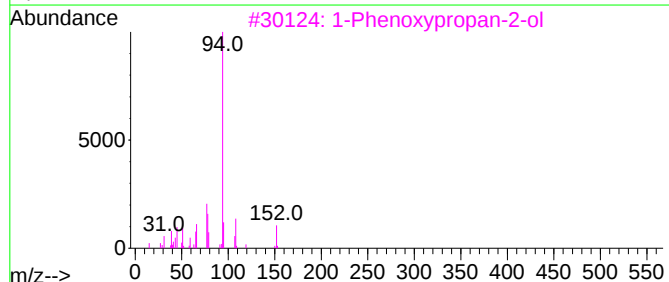
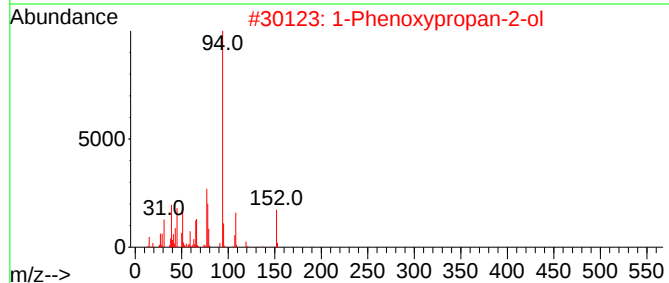
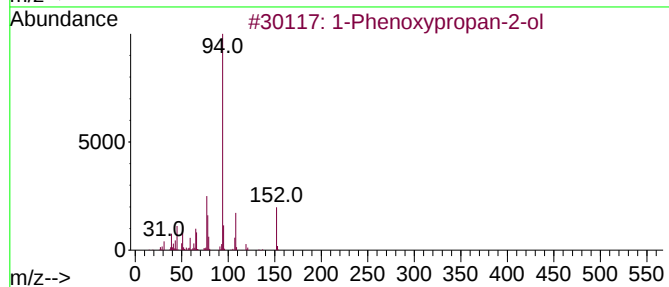
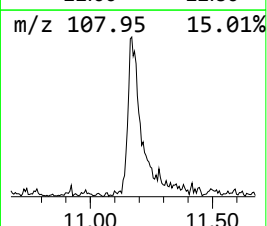
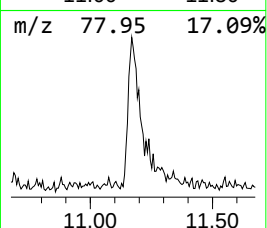
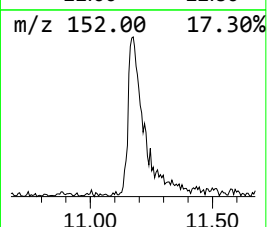
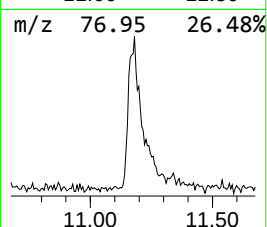
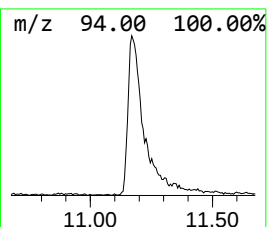
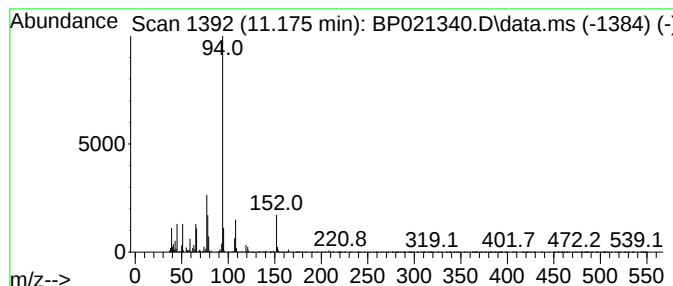
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 1-Phenoxypropan-2-ol Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.175	2.51 ng/ul	185231	Naphthalene-d8	10.381

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	96
2			1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	96
3			1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	95
4			1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	93
5			1-Propanol, 3-phenoxy-	152	C9H12O2	006180-61-6	87



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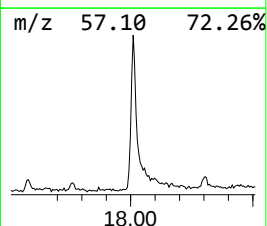
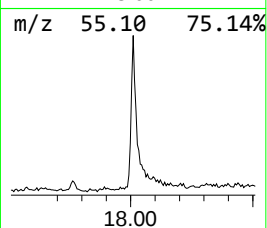
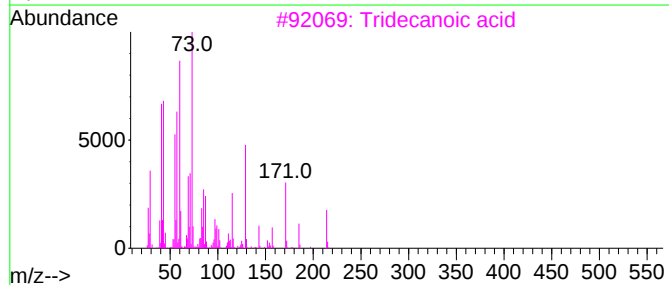
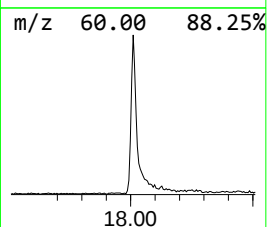
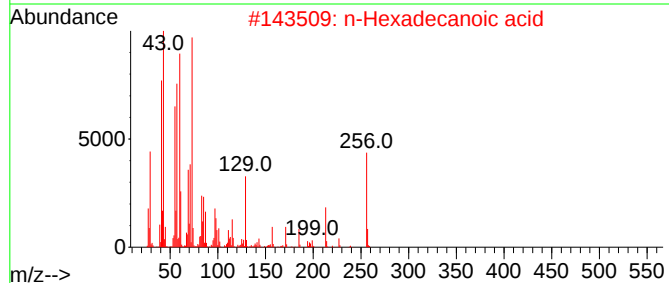
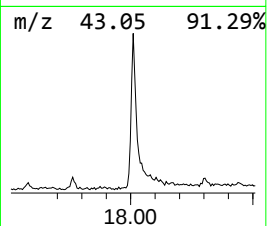
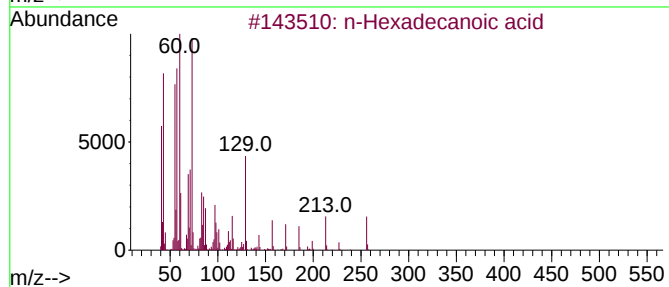
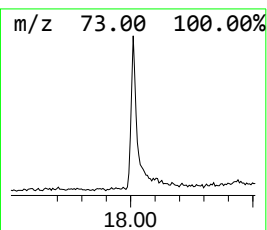
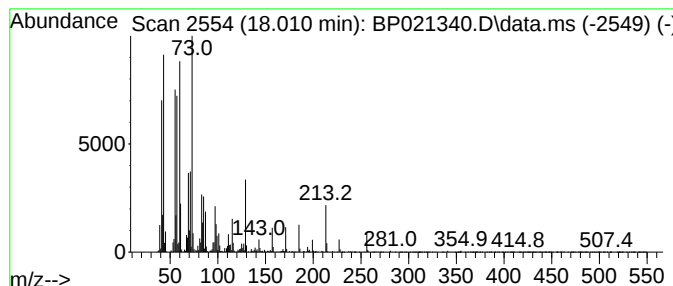
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TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 n-Hexadecanoic acid Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.010	4.17 ng/ul	546594	Phenanthrene-d10	17.081

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	Tridecanoic acid	214	C13H26O2	000638-53-9	96		
4	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	95		
5	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	93		



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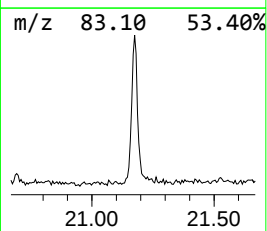
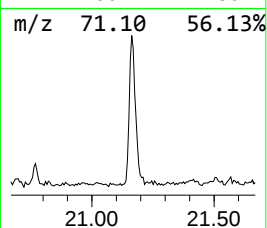
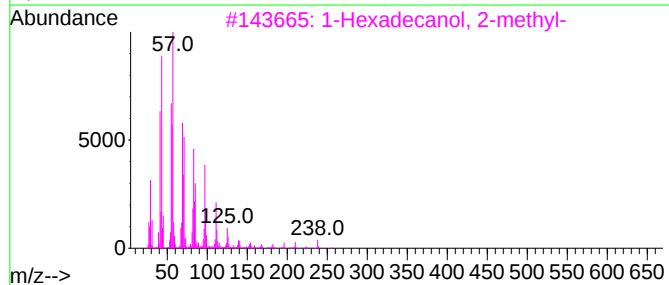
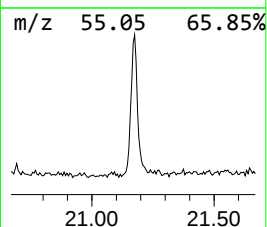
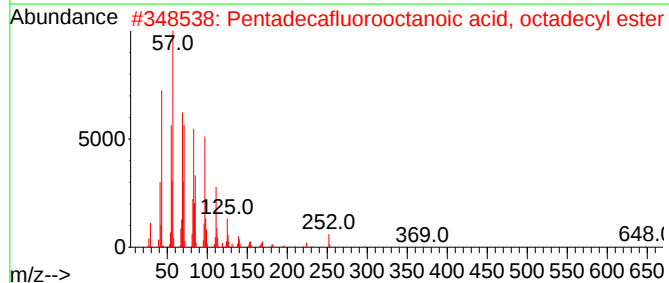
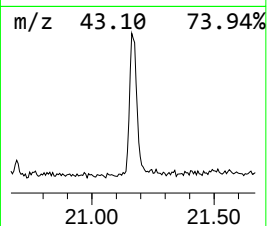
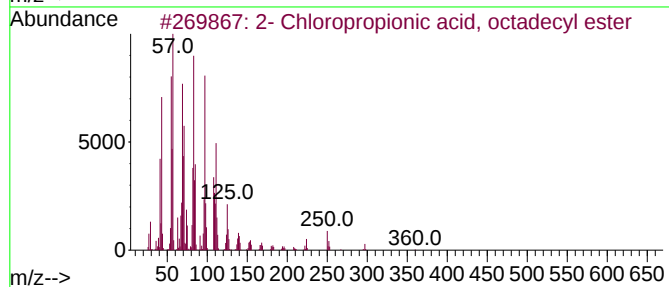
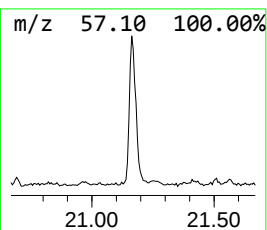
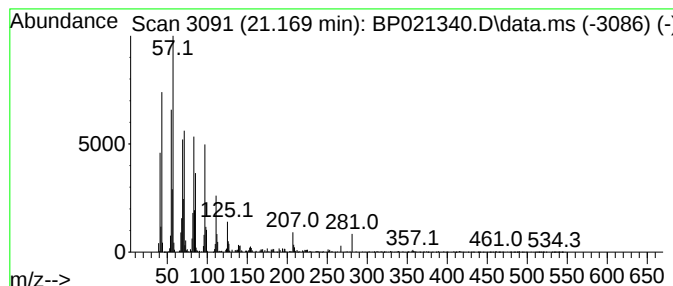
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Quant Title : SVOA CALIBRATION

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

Peak Number 3 2- Chloropropionic acid, oc... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.169	4.01 ng/ul	587172	Chrysene-d12	21.528

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-	Chloropropionic acid, octadec...	360	C21H41ClO2	088104-31-8	94	
2	Pentadecafluorooctanoic acid, oc...	666	C26H37F15O2	1000406-04-8	93		
3	1-Hexadecanol, 2-methyl-	256	C17H36O	002490-48-4	93		
4	Pentadecafluorooctanoic acid, he...	638	C24H33F15O2	1000406-04-6	93		
5	Cyclotriacontane	420	C30H60	000297-35-8	93		



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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
1-Phenoxypropan...	11.175	2.5	ng/ul	185231	2	10.381	1478720	20.0
n-Hexadecanoic ...	18.010	4.2	ng/ul	546594	4	17.081	2623180	20.0
2- Chloropropio...	21.169	4.0	ng/ul	587172	5	21.528	2925500	20.0