

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102121\
 Data File : BP007595.D
 Acq On : 22 Oct 2021 18:05
 Operator : CG/JU
 Sample : M4281-16
 Misc :
 ALS Vial : 55 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BFY75

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-BP101421.M
 Title : SVOA CALIBRATION

Signal : TIC: BP007595.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.375	11	15	22	rVB	40272	51405	1.01%	0.110%
2	3.652	60	62	68	rBV	11872	15457	0.30%	0.033%
3	3.793	83	86	102	rBV	180428	299833	5.89%	0.641%
4	4.034	122	127	133	rVB	25080	36563	0.72%	0.078%
5	4.128	139	143	148	rVB	13142	19332	0.38%	0.041%
6	4.222	155	159	165	rVB8	10918	17847	0.35%	0.038%
7	4.581	218	220	225	rVB2	6052	7097	0.14%	0.015%
8	5.052	297	300	304	rVB5	7018	9013	0.18%	0.019%
9	5.646	396	401	405	rVB8	3573	7393	0.15%	0.016%
10	6.028	461	466	470	rBV3	6153	8413	0.17%	0.018%
11	6.328	513	517	522	rVB5	10045	14119	0.28%	0.030%
12	7.105	643	649	662	rVB	183283	318854	6.27%	0.681%
13	7.275	669	678	690	rBV	529950	855682	16.82%	1.829%
14	7.393	695	698	703	rBV6	2392	5200	0.10%	0.011%
15	7.475	703	712	720	rBV	614149	950957	18.69%	2.032%
16	7.552	722	725	729	rBV4	4252	6697	0.13%	0.014%
17	7.946	785	792	799	rBV	705247	1096762	21.56%	2.344%
18	8.016	801	804	812	rVB5	8470	13280	0.26%	0.028%
19	8.193	826	834	840	rBV5	6097	13645	0.27%	0.029%
20	8.652	904	912	922	rBV	514539	819437	16.11%	1.751%
21	9.104	981	989	1000	rBV	641942	1087387	21.37%	2.324%
22	9.240	1005	1012	1023	rVV	86459	151928	2.99%	0.325%
23	9.446	1042	1047	1051	rVB8	3916	5262	0.10%	0.011%
24	9.569	1062	1068	1078	rVB4	14647	29653	0.58%	0.063%
25	9.828	1103	1112	1121	rBV	524742	873103	17.16%	1.866%
26	10.046	1146	1149	1155	rBV8	2608	5222	0.10%	0.011%
27	10.369	1196	1204	1215	rBV	894957	1476629	29.02%	3.155%
28	10.534	1226	1232	1241	rBV	205526	330380	6.49%	0.706%
29	10.751	1259	1269	1281	rVB	932576	1578531	31.03%	3.373%
30	10.881	1283	1291	1299	rBV	742766	1254078	24.65%	2.680%
31	11.169	1334	1340	1343	rVB8	6981	11315	0.22%	0.024%
32	11.398	1374	1379	1383	rVB8	3747	6340	0.12%	0.014%
33	12.063	1489	1492	1498	rVB4	11356	18974	0.37%	0.041%
34	12.340	1531	1539	1549	rBV2	19705	41577	0.82%	0.089%
35	12.793	1611	1616	1620	rBV7	4034	8475	0.17%	0.018%
36	12.957	1639	1644	1647	rBV7	5167	7246	0.14%	0.015%

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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-BP101421.M
 Title : SVOA CALIBRATION

37	13.075	1660	1664	1667	rBV6	3745	6198	0.12%	0.013%
38	13.204	1677	1686	1691	rVB10	9362	16124	0.32%	0.034%
39	13.316	1700	1705	1708	rBV6	3708	6058	0.12%	0.013%
40	13.428	1718	1724	1726	rBV6	5651	8468	0.17%	0.018%
41	13.492	1730	1735	1741	rVV10	6818	13865	0.27%	0.030%
42	13.557	1741	1746	1752	rVB9	6478	8876	0.17%	0.019%
43	13.904	1798	1805	1806	rBV7	2886	5280	0.10%	0.011%
44	13.992	1813	1820	1829	rBV	1623311	2258284	44.39%	4.826%
45	14.128	1841	1843	1849	rVB6	3955	5684	0.11%	0.012%
46	14.275	1857	1868	1879	rBV	1818022	2633498	51.76%	5.628%
47	14.428	1890	1894	1896	rBV5	3082	5440	0.11%	0.012%
48	14.498	1902	1906	1912	rVB2	16278	24169	0.48%	0.052%
49	14.581	1913	1920	1934	rVV	1463185	2172480	42.70%	4.642%
50	14.769	1943	1952	1957	rBV	186759	278643	5.48%	0.595%
51	14.998	1989	1991	1994	rVB4	5386	5608	0.11%	0.012%
52	15.134	2008	2014	2027	rBV2	61178	183901	3.61%	0.393%
53	15.234	2027	2031	2035	rVB4	17689	27841	0.55%	0.059%
54	15.398	2055	2059	2064	rVB4	12647	19679	0.39%	0.042%
55	15.575	2082	2089	2100	rBV	2366862	3446443	67.74%	7.365%
56	15.686	2102	2108	2116	rVV	131510	192753	3.79%	0.412%
57	15.816	2125	2130	2136	rVB2	29487	43186	0.85%	0.092%
58	16.022	2162	2165	2170	rVB7	10908	15059	0.30%	0.032%
59	16.134	2181	2184	2185	rBV3	7843	9347	0.18%	0.020%
60	16.263	2198	2206	2211	rBV3	9611	27964	0.55%	0.060%
61	16.310	2212	2214	2218	rVB5	5683	5356	0.11%	0.011%
62	16.363	2218	2223	2228	rBV8	10197	16395	0.32%	0.035%
63	16.733	2283	2286	2291	rBV7	8568	13762	0.27%	0.029%
64	16.986	2325	2329	2341	rBV	24978	58934	1.16%	0.126%
65	17.151	2355	2357	2361	rVB5	9160	9753	0.19%	0.021%
66	17.333	2381	2388	2396	rBV	1819573	2633677	51.76%	5.628%
67	17.433	2398	2405	2417	rVV	2936840	4047871	79.56%	8.650%
68	17.575	2424	2429	2435	rVB2	30999	52745	1.04%	0.113%
69	18.016	2500	2504	2511	rVB2	41370	48064	0.94%	0.103%
70	18.228	2536	2540	2547	rBV	141290	191692	3.77%	0.410%
71	18.598	2601	2603	2608	rVB6	14400	13057	0.26%	0.028%
72	18.951	2660	2663	2667	rBV2	29044	43671	0.86%	0.093%
73	19.245	2711	2713	2717	rVB	36103	39157	0.77%	0.084%
74	19.316	2721	2725	2728	rBV	45808	61870	1.22%	0.132%
75	19.669	2779	2785	2791	rBV	3617801	4976544	97.81%	10.634%
76	21.133	3031	3034	3042	rBV	256153	352424	6.93%	0.753%

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Integrator: RTE
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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

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Title : SVOA CALIBRATION

77	21.421	3078	3083	3089	rBV	2301929	3036916	59.69%	6.490%
78	23.710	3465	3472	3482	rVB2	2551420	5087919	100.00%	10.872%
79	23.868	3492	3499	3511	rVB2	1529537	3238679	63.65%	6.921%

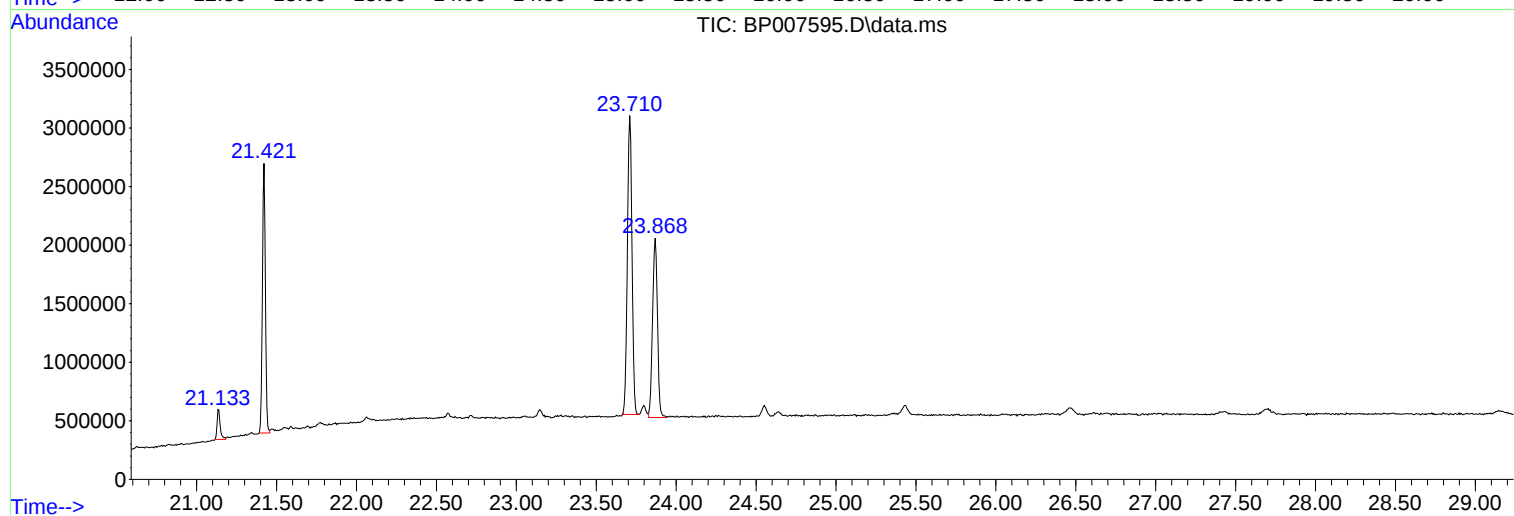
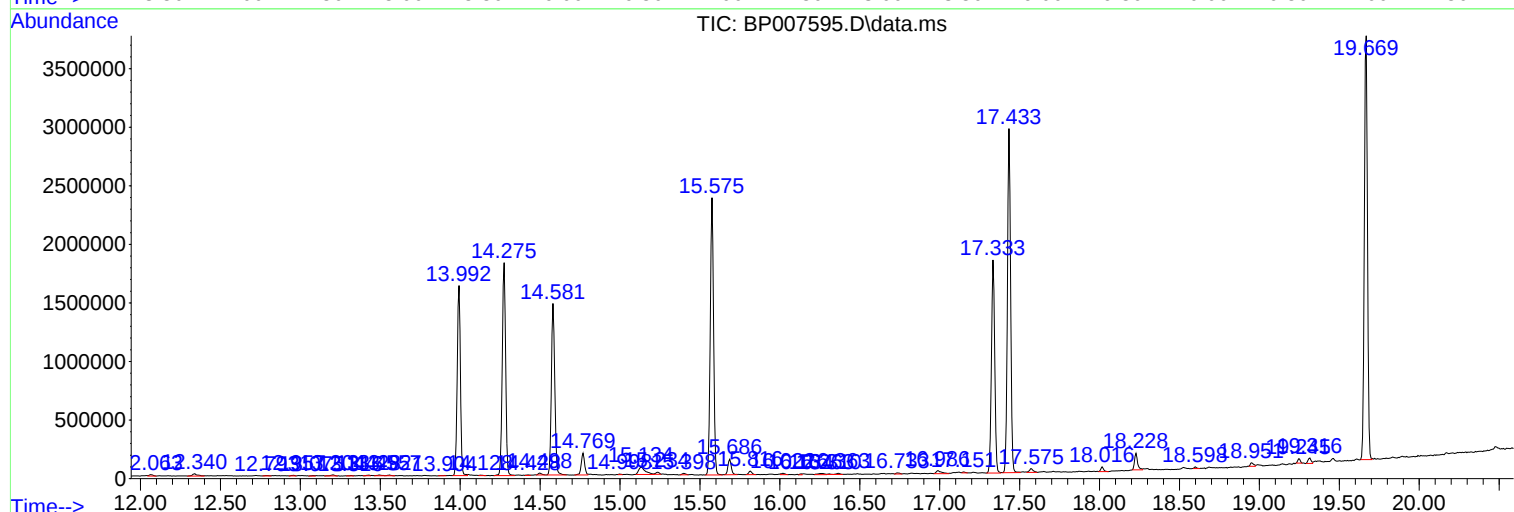
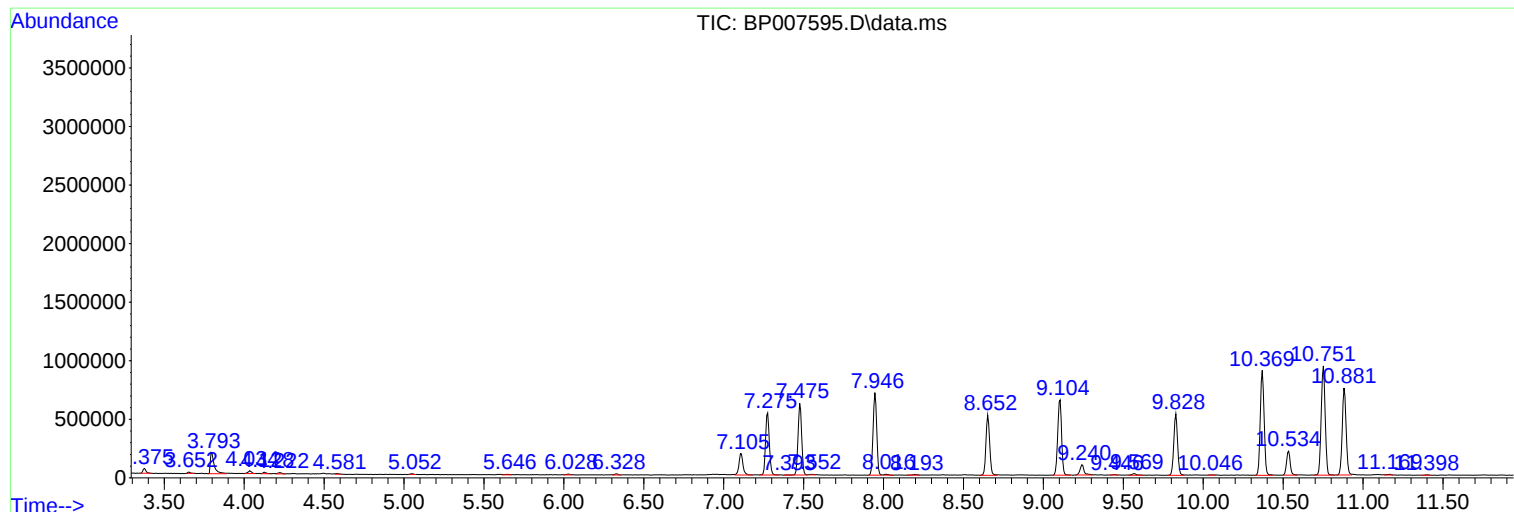
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP101421.M
Quant Title : SVOA CALIBRATION

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



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Sample : M4281-16
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Instrument :
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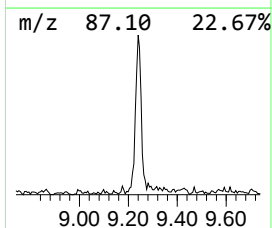
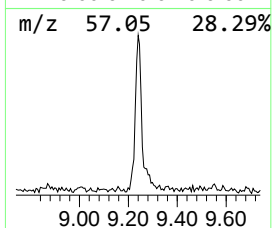
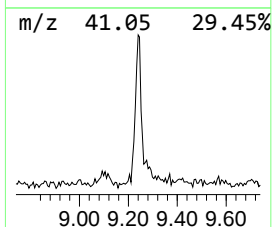
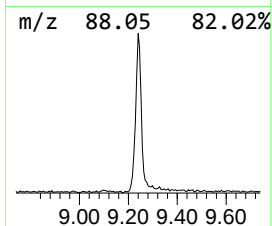
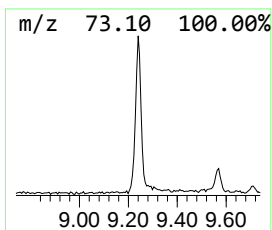
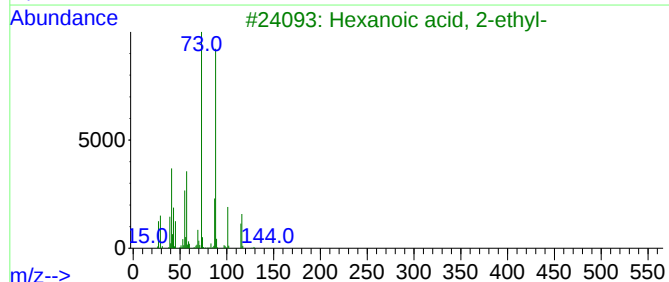
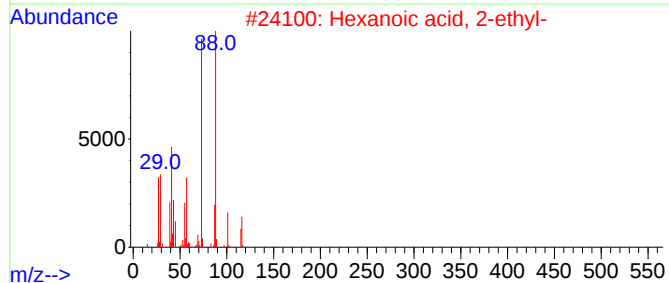
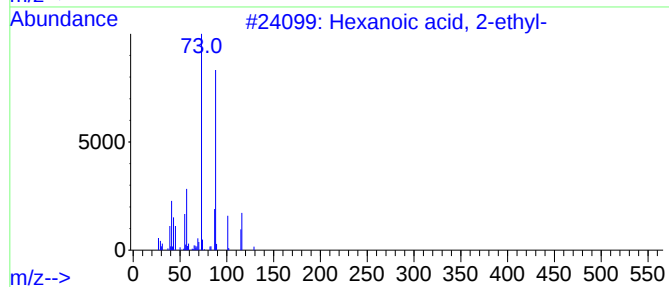
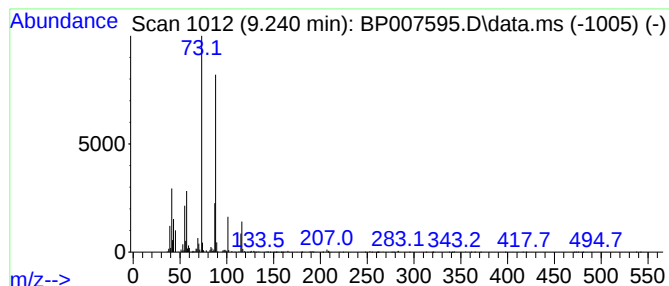
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP101421.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Hexanoic acid, 2-ethyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.240	2.77 ng/ul	151928	1,4-Dichlorobenzene-d4	7.946

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexanoic acid, 2-ethyl-	144	C8H16O2	000149-57-5	90
2			Hexanoic acid, 2-ethyl-	144	C8H16O2	000149-57-5	90
3			Hexanoic acid, 2-ethyl-	144	C8H16O2	000149-57-5	90
4			Hexanoic acid, 2-ethyl-	144	C8H16O2	000149-57-5	78
5			Hexanoic acid, 2-ethyl-	144	C8H16O2	000149-57-5	76



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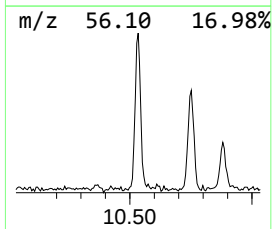
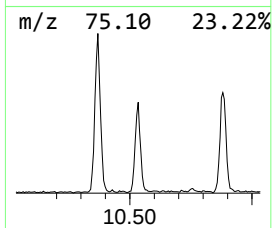
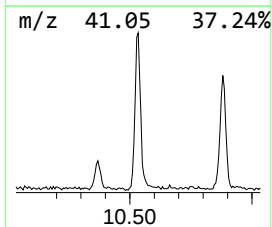
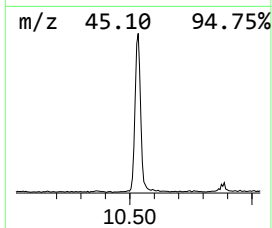
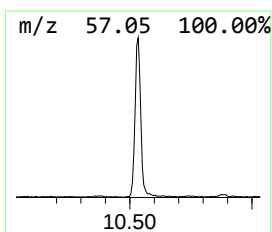
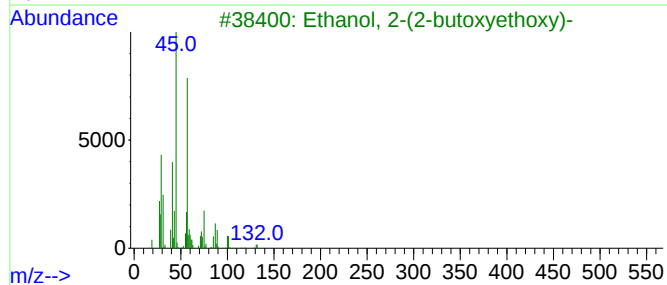
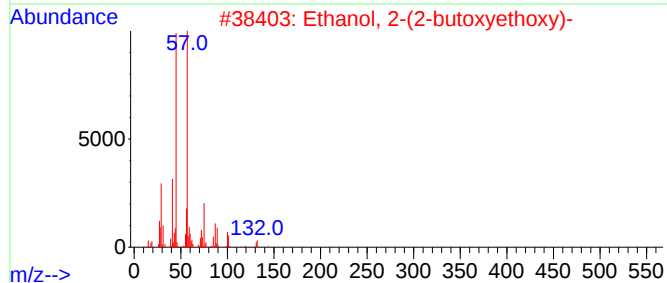
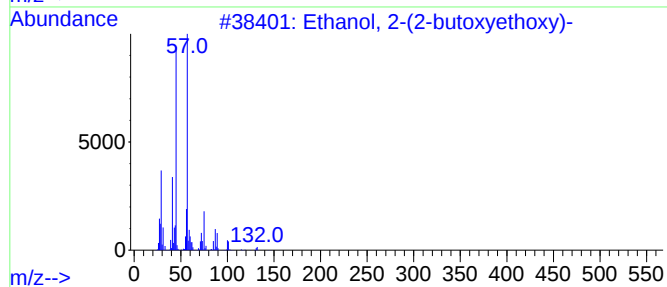
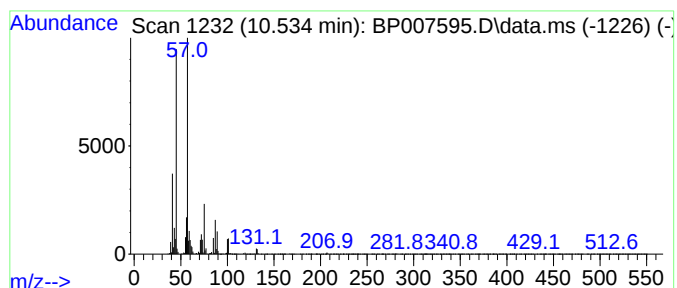
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP101421.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 Ethanol, 2-(2-butoxyethoxy)- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.534	4.19 ng/ul	330380	Naphthalene-d8	10.752

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	90
2			Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	90
3			Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	83
4			Ethanol, 1-(2-butoxyethoxy)-	162	C8H18O3	054446-78-5	83
5			Ethanol, 2-[2-(2-methylpropoxy)e...	162	C8H18O3	018912-80-6	64



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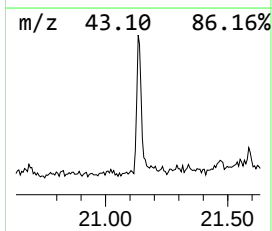
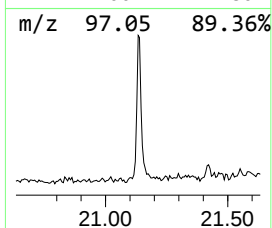
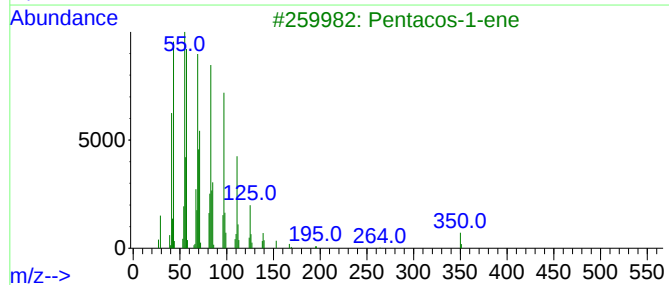
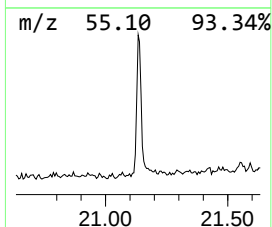
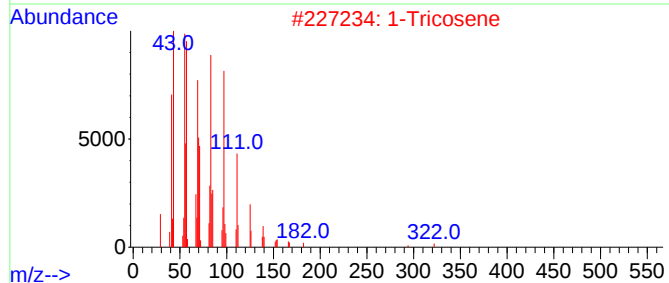
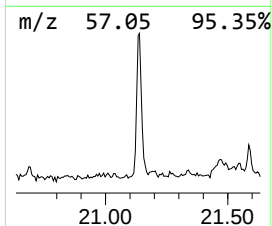
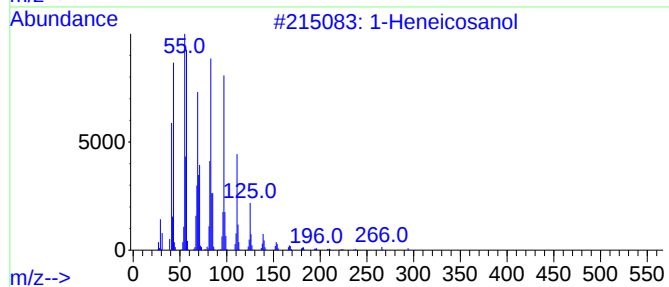
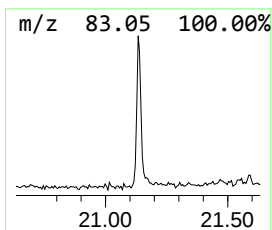
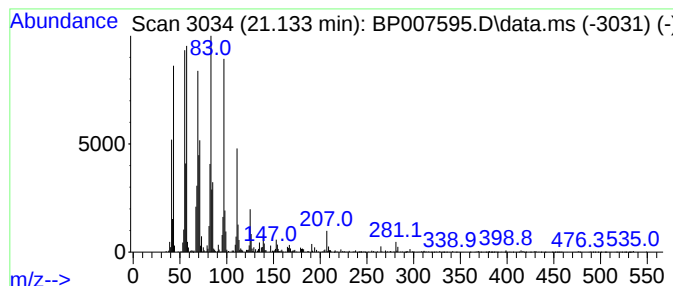
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP101421.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 1-Heneicosanol Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.133	2.32 ng/ul	352424	Chrysene-d12	21.422

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Heneicosanol	312	C21H44O	015594-90-8	95
2			1-Tricosene	322	C23H46	018835-32-0	95
3			Pentacos-1-ene	350	C25H50	016980-85-1	94
4			n-Nonadecanol-1	284	C19H40O	001454-84-8	94
5			Pentafluoropropionic acid, penta...	374	C18H31F5O2	959092-08-1	94



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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
Hexanoic acid, ...	9.240	2.8	ng/ul	151928	1	7.946	1096760	20.0
Ethanol, 2-(2-b...	10.534	4.2	ng/ul	330380	2	10.752	1578530	20.0
1-Heneicosanol	21.133	2.3	ng/ul	352424	5	21.422	3036920	20.0