

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP081021\
 Data File : BP006629.D
 Acq On : 10 Aug 2021 18:25
 Operator : CG/JU
 Sample : M3208-06
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C00E6

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-BP081021.M
 Title : SVOA CALIBRATION

Signal : TIC: BP006629.D

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.640	91	94	106	rVB	185435	291405	10.64%	0.551%
2	6.899	643	648	655	rVB	528010	796033	29.06%	1.506%
3	7.064	671	676	691	rVB	446289	732793	26.75%	1.387%
4	7.252	703	708	718	rVB	545622	849473	31.01%	1.607%
5	7.722	780	788	796	rVB	620935	1048511	38.28%	1.984%
6	8.434	903	909	916	rVB	486858	771321	28.16%	1.460%
7	8.875	978	984	991	rVV	546134	869139	31.73%	1.645%
8	8.946	994	996	1001	rVB3	18560	23962	0.87%	0.045%
9	9.593	1100	1106	1120	rBV	453214	728718	26.60%	1.379%
10	10.128	1190	1197	1207	rVB	618399	1020800	37.27%	1.932%
11	10.299	1220	1226	1236	rBV	488491	813066	29.68%	1.539%
12	10.505	1253	1261	1267	rBV	901116	1493709	54.53%	2.827%
13	10.552	1267	1269	1277	rVB2	23663	34621	1.26%	0.066%
14	10.652	1279	1286	1299	rBV	289524	543328	19.83%	1.028%
15	11.540	1432	1437	1445	rVB3	19999	35100	1.28%	0.066%
16	11.940	1501	1505	1512	rVB	30630	44043	1.61%	0.083%
17	12.099	1525	1532	1536	rVV5	13244	28711	1.05%	0.054%
18	12.169	1538	1544	1551	rVB	38131	61221	2.23%	0.116%
19	12.387	1575	1581	1589	rVB	20365	34160	1.25%	0.065%
20	12.634	1617	1623	1629	rBV9	8315	18385	0.67%	0.035%
21	12.904	1663	1669	1673	rVV	15275	26344	0.96%	0.050%
22	13.193	1713	1718	1724	rVB2	49820	68852	2.51%	0.130%
23	13.522	1769	1774	1775	rBV4	14295	20446	0.75%	0.039%
24	13.681	1795	1801	1805	rVV	25636	44059	1.61%	0.083%
25	13.734	1805	1810	1812	rVV2	19000	28440	1.04%	0.054%
26	13.781	1812	1818	1826	rVV	1133505	1530186	55.86%	2.896%
27	13.851	1826	1830	1834	rVV5	16205	24864	0.91%	0.047%
28	13.910	1834	1840	1844	rVB5	12000	23577	0.86%	0.045%
29	14.051	1853	1864	1875	rBV	1209868	1836721	67.05%	3.476%
30	14.269	1897	1901	1907	rVB	44876	60344	2.20%	0.114%
31	14.357	1907	1916	1923	rBV	1306296	1874124	68.42%	3.546%
32	14.569	1946	1952	1965	rVV	525730	739242	26.99%	1.399%
33	14.728	1974	1979	1981	rBV2	26213	40753	1.49%	0.077%
34	14.757	1981	1984	1988	rVV2	25491	44288	1.62%	0.084%
35	14.834	1993	1997	2004	rVV3	18039	27783	1.01%	0.053%
36	14.981	2014	2022	2027	rBV5	13284	28454	1.04%	0.054%

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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-BP081021.M
 Title : SVOA CALIBRATION

37	15.034	2027	2031	2035	rVB2	16217	21879	0.80%	0.041%
38	15.228	2060	2064	2068	rVB	48167	55678	2.03%	0.105%
39	15.357	2079	2086	2092	rBV	1508425	2162678	78.95%	4.092%
40	15.481	2099	2107	2116	rBV	482025	650253	23.74%	1.230%
41	15.593	2122	2126	2130	rBV	15982	22905	0.84%	0.043%
42	15.640	2130	2134	2139	rVB5	12603	20563	0.75%	0.039%
43	15.704	2139	2145	2152	rBV3	20169	39233	1.43%	0.074%
44	15.851	2167	2170	2178	rVB5	15997	33194	1.21%	0.063%
45	15.945	2178	2186	2192	rBV7	14005	29253	1.07%	0.055%
46	16.093	2207	2211	2218	rBV	63513	95127	3.47%	0.180%
47	16.163	2218	2223	2231	rVB9	9460	24218	0.88%	0.046%
48	16.663	2301	2308	2310	rBV5	15290	32696	1.19%	0.062%
49	16.769	2322	2326	2334	rVB2	29397	55466	2.02%	0.105%
50	16.845	2334	2339	2343	rBV3	19981	39032	1.42%	0.074%
51	16.892	2343	2347	2351	rVV2	61141	79130	2.89%	0.150%
52	16.934	2351	2354	2362	rVB6	18148	30908	1.13%	0.058%
53	17.028	2366	2370	2377	rBV4	36496	54380	1.99%	0.103%
54	17.110	2377	2384	2389	rBV	1391641	2139637	78.11%	4.049%
55	17.157	2389	2392	2396	rVV	148156	191958	7.01%	0.363%
56	17.210	2396	2401	2412	rVB2	1509942	2157196	78.75%	4.082%
57	17.310	2412	2418	2422	rBV8	14545	29223	1.07%	0.055%
58	17.428	2434	2438	2443	rVB2	29629	42612	1.56%	0.081%
59	17.634	2469	2473	2476	rBV	58902	71205	2.60%	0.135%
60	17.669	2476	2479	2481	rVV3	14540	21162	0.77%	0.040%
61	17.698	2482	2484	2488	rVB3	19511	20291	0.74%	0.038%
62	17.804	2498	2502	2510	rVB2	68070	89193	3.26%	0.169%
63	17.904	2515	2519	2524	rVV7	29573	42537	1.55%	0.080%
64	17.963	2525	2529	2530	rVV	35733	39543	1.44%	0.075%
65	17.987	2530	2533	2535	rVV2	52040	68039	2.48%	0.129%
66	18.022	2535	2539	2547	rVV2	335371	540623	19.74%	1.023%
67	18.092	2547	2551	2557	rVV	79739	134263	4.90%	0.254%
68	18.169	2560	2564	2568	rVV2	61928	101397	3.70%	0.192%
69	18.210	2568	2571	2576	rVB2	42701	54895	2.00%	0.104%
70	18.310	2583	2588	2594	rBV2	71539	113142	4.13%	0.214%
71	18.475	2612	2616	2620	rBV	62102	83181	3.04%	0.157%
72	18.763	2662	2665	2668	rBV3	17870	20133	0.73%	0.038%
73	18.798	2668	2671	2675	rVB3	16450	18862	0.69%	0.036%
74	18.881	2680	2685	2689	rBV	55959	88915	3.25%	0.168%
75	18.922	2689	2692	2695	rBV2	84666	105599	3.85%	0.200%
76	19.028	2704	2710	2716	rBV9	21977	57385	2.09%	0.109%

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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-BP081021.M
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77	19.134	2724	2728	2732	rVB	208643	265122	9.68%	0.502%
78	19.198	2737	2739	2742	rVB	20474	21211	0.77%	0.040%
79	19.263	2747	2750	2757	rVB3	76908	108006	3.94%	0.204%
80	19.457	2778	2783	2796	rBV	1878752	2739298	100.00%	5.184%
81	19.563	2797	2801	2806	rVB4	31300	47575	1.74%	0.090%
82	19.716	2824	2827	2831	rBV6	23312	29388	1.07%	0.056%
83	19.975	2868	2871	2873	rBV4	64725	85890	3.14%	0.163%
84	20.004	2873	2876	2881	rVB	190993	212369	7.75%	0.402%
85	20.492	2955	2959	2963	rVB	130465	144645	5.28%	0.274%
86	20.945	3032	3036	3043	rBV	1374557	1699958	62.06%	3.217%
87	21.216	3076	3082	3086	rBV	1807468	2505527	91.47%	4.741%
88	21.375	3106	3109	3114	rBV	202032	254316	9.28%	0.481%
89	21.563	3138	3141	3150	rVB2	225648	248859	9.08%	0.471%
90	21.828	3182	3186	3187	rBV	649666	724392	26.44%	1.371%
91	21.851	3187	3190	3200	rVB	987517	1441115	52.61%	2.727%
92	22.310	3265	3268	3271	rVB	191245	214052	7.81%	0.405%
93	22.551	3304	3309	3319	rBV	1308470	1794107	65.50%	3.395%
94	22.845	3354	3359	3363	rBV	953333	1440608	52.59%	2.726%
95	22.898	3363	3368	3378	rVB	1361478	2487064	90.79%	4.706%
96	23.380	3444	3450	3456	rBV2	1106000	2033494	74.23%	3.848%
97	23.527	3469	3475	3484	rVB2	1299791	2551696	93.15%	4.829%
98	23.786	3513	3519	3530	rVB	1190173	2214475	80.84%	4.190%
99	24.151	3575	3581	3589	rVB	728255	1496999	54.65%	2.833%
100	24.245	3591	3597	3608	rBV	725806	1750511	63.90%	3.313%

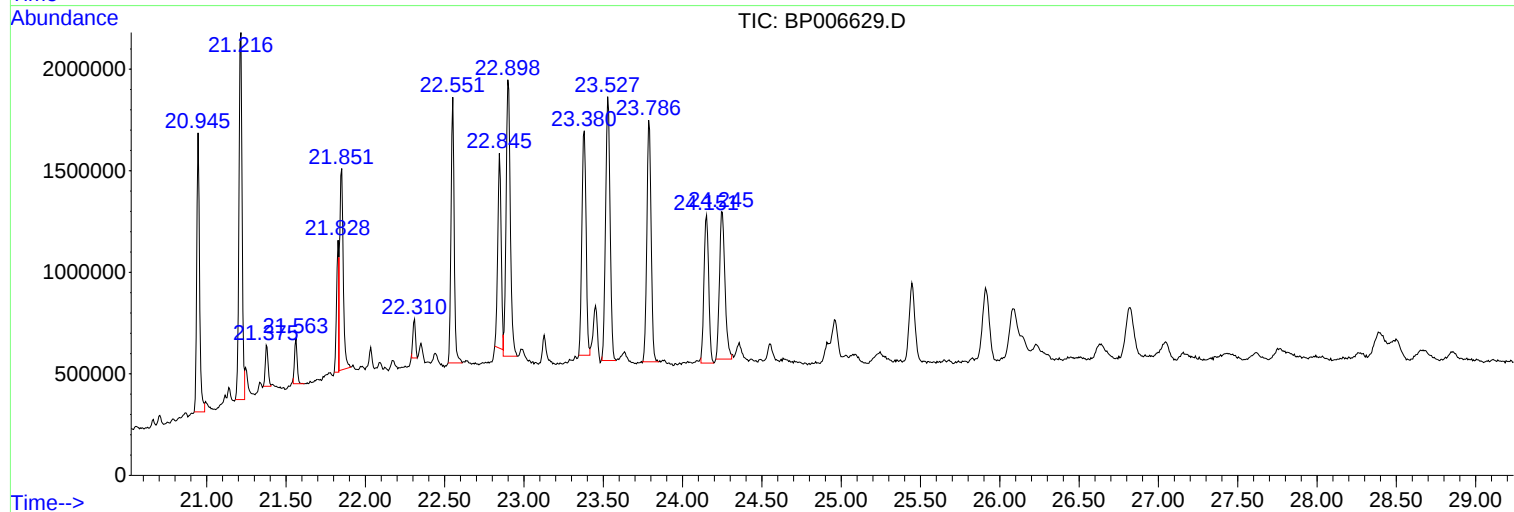
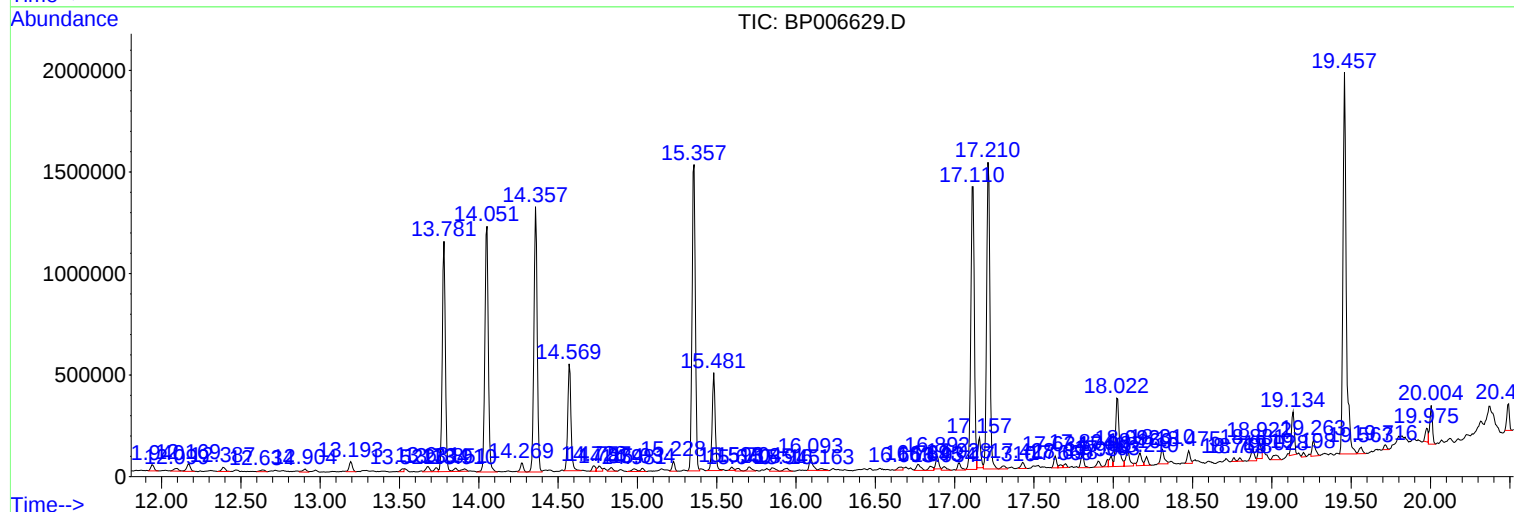
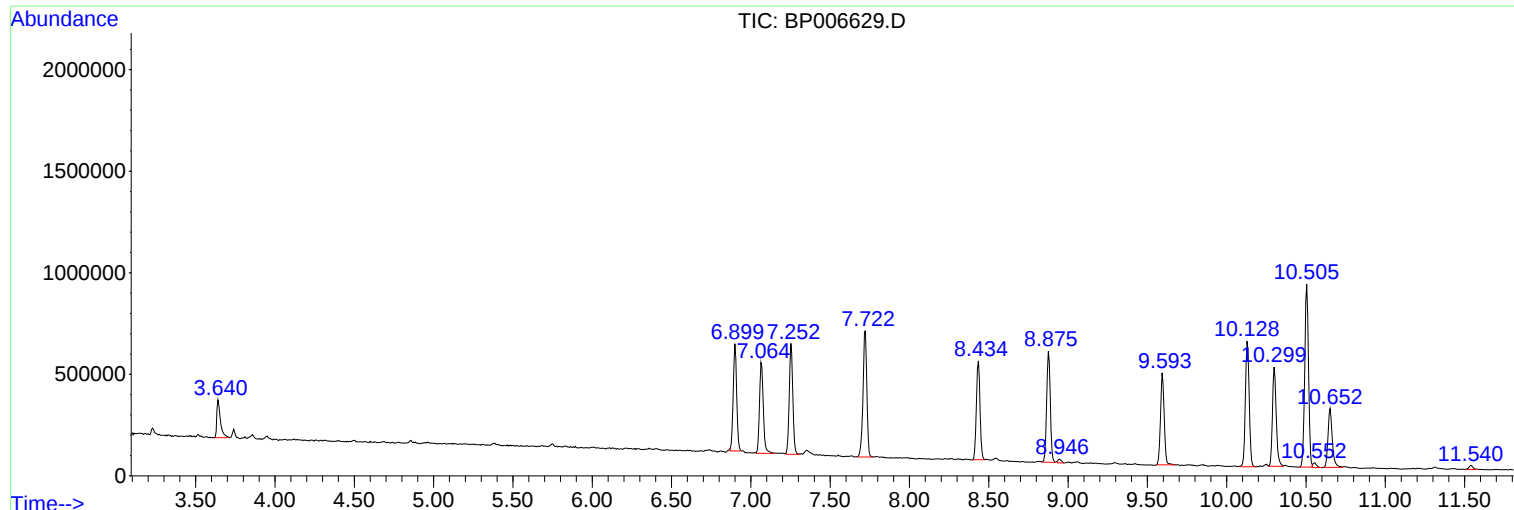
Sum of corrected areas: 52845237

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ALS Vial : 14 Sample Multiplier: 1

Instrument :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP081021.M
Quant Title : SVOA CALIBRATION

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



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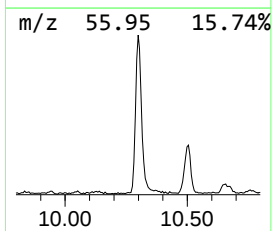
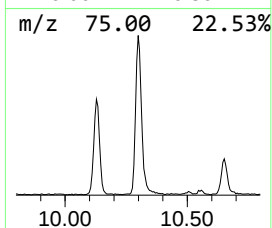
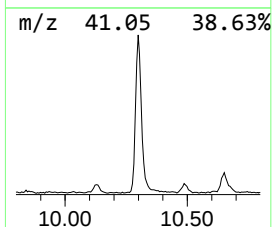
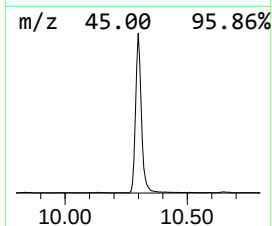
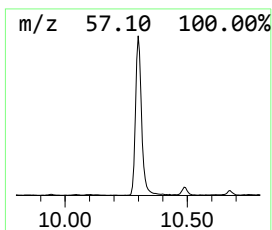
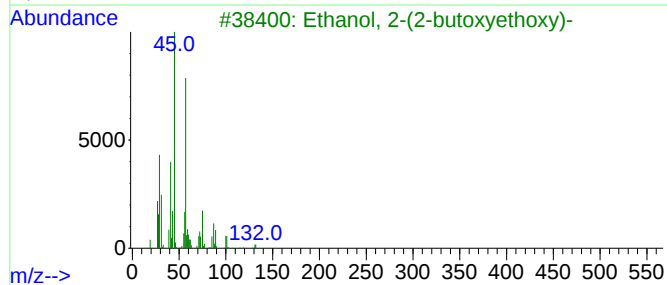
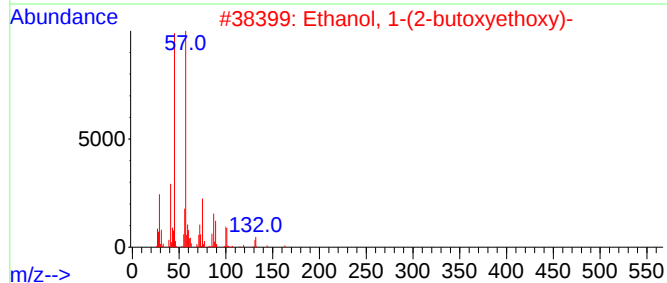
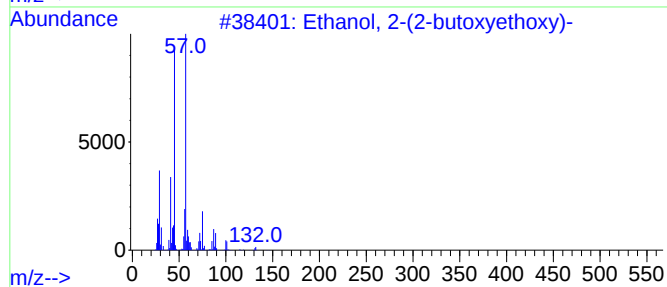
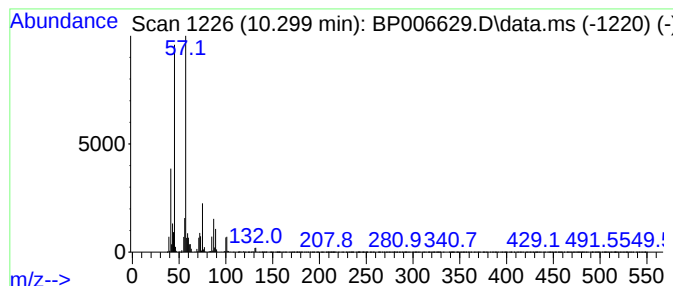
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP081021.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Ethanol, 2-(2-butoxyethoxy)- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.300	10.89 ng/ul	813066	Naphthalene-d8	10.505

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	90
4			Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	83
5			Ethylene glycol monoisobutyl ether	118	C6H14O2	004439-24-1	64



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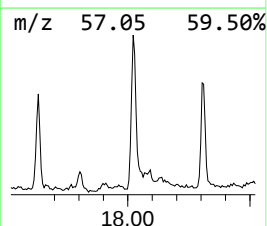
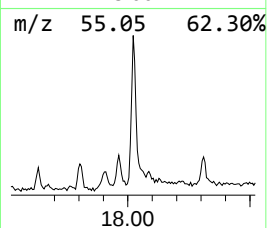
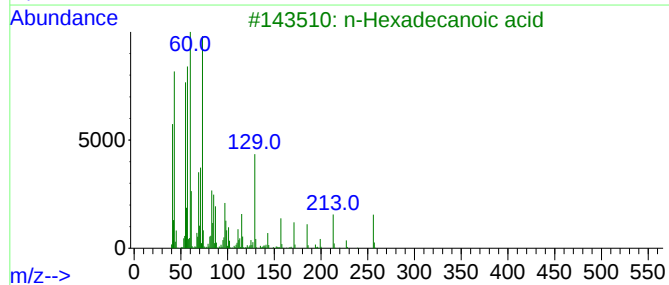
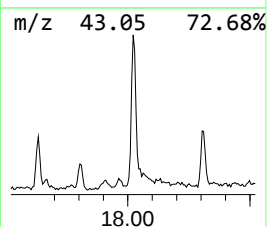
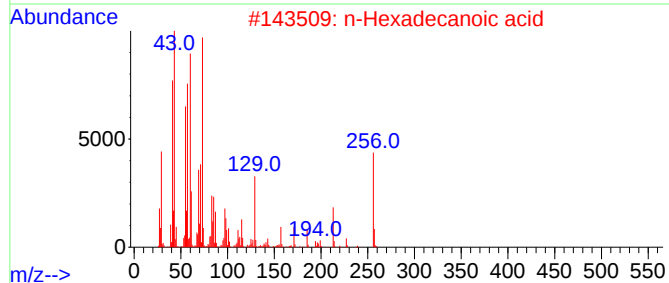
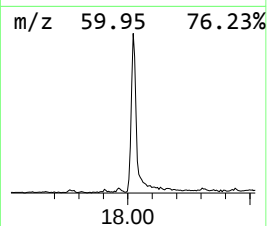
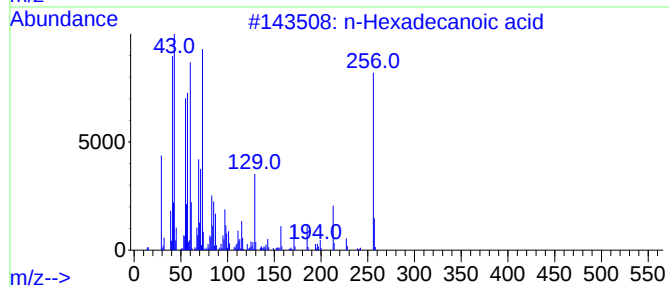
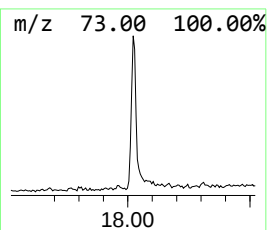
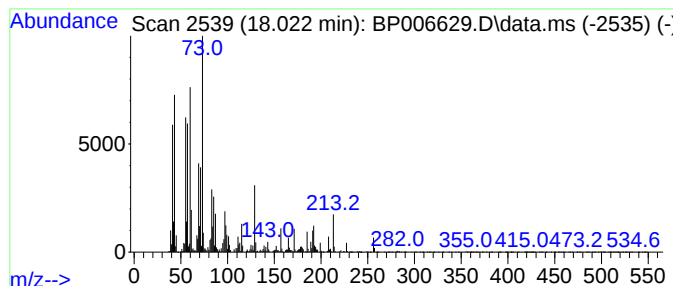
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP081021.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 n-Hexadecanoic acid Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.020	5.05 ng/ul	540623	Phenanthrene-d10	17.116

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	96
4			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	95
5			Octadecanoic acid	284	C18H36O2	000057-11-4	81



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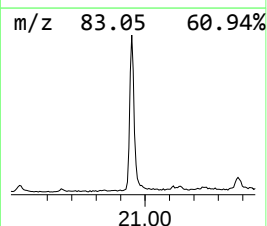
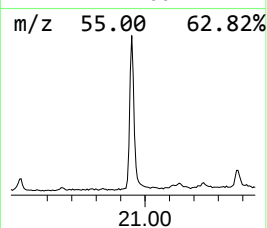
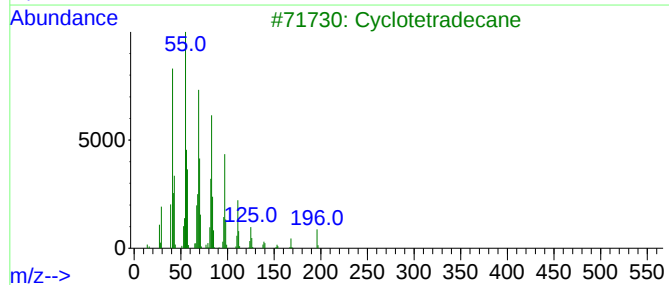
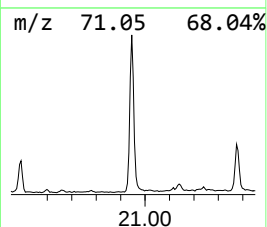
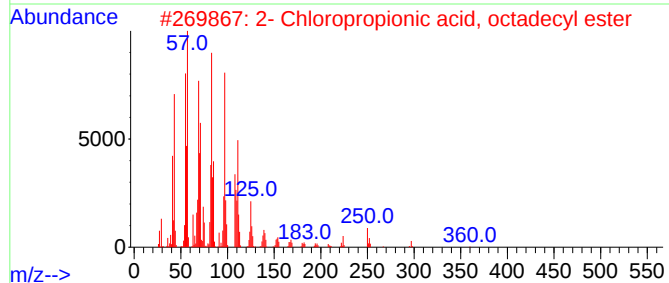
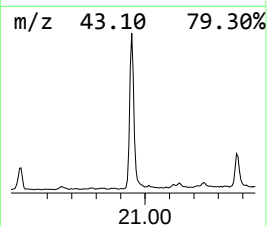
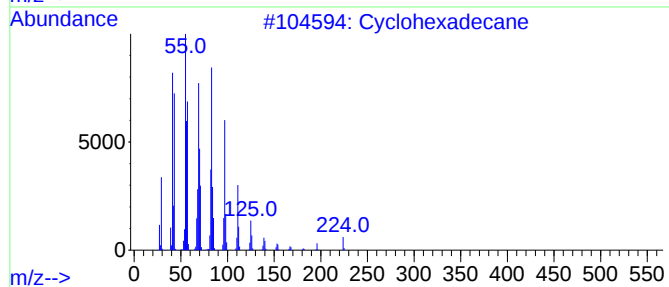
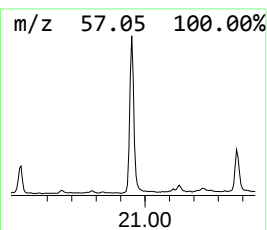
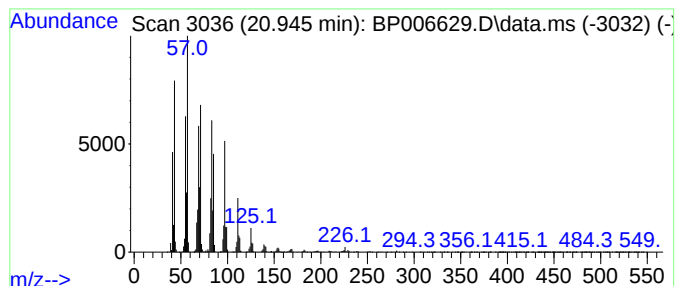
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Quant Title : SVOA CALIBRATION

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

Peak Number 3 (DEL) Alkane: Cyclic20.95 Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.950	13.57 ng/ul	1699960	Chrysene-d12	21.216

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexadecane	224	C16H32	000295-65-8	93
2			2- Chloropropionic acid, octadec...	360	C21H41ClO2	088104-31-8	93
3			Cyclotetradecane	196	C14H28	000295-17-0	91
4			Pentadecafluorooctanoic acid, oc...	666	C26H37F15O2	1000406-04-8	90
5			Pentafluoropropionic acid, octad...	416	C21H37F5O2	959261-25-7	81



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP081021\
Data File : BP006629.D
Acq On : 10 Aug 2021 18:25
Operator : CG/JU
Sample : M3208-06
Misc :
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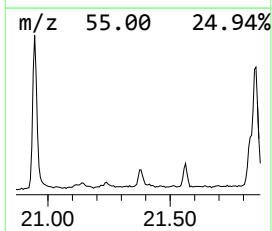
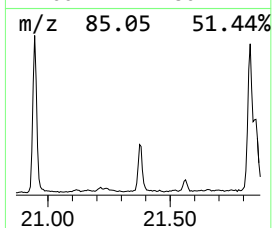
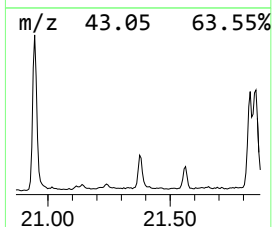
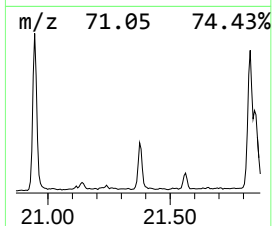
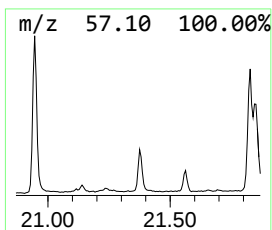
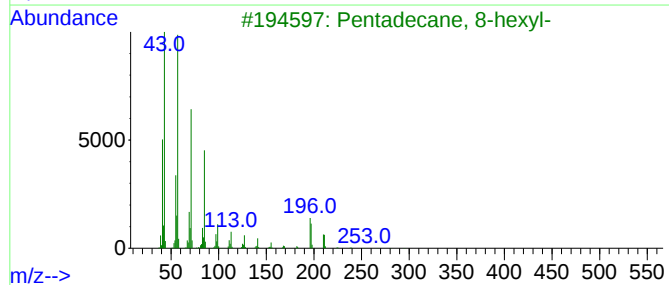
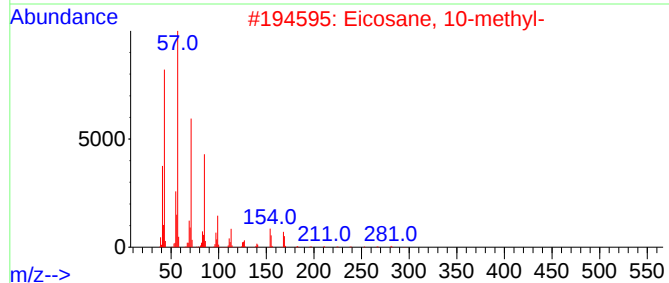
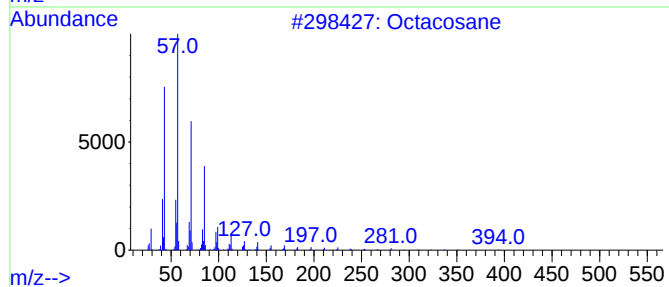
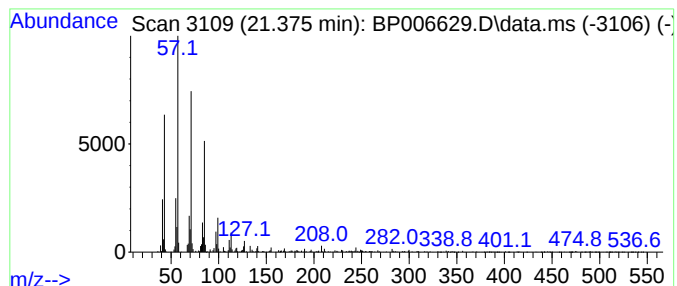
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 (DEL) Alkane: Straight-Chai... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.	
21.370	2.03 ng/ul	254316	Chrysene-d12	21.216	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octacosane	394	C28H58	000630-02-4	94
2	Eicosane, 10-methyl-	296	C21H44	054833-23-7	94
3	Pentadecane, 8-hexyl-	296	C21H44	013475-75-7	94
4	Tridecane, 1-iodo-	310	C13H27I	035599-77-0	93
5	Docosane, 11-butyl-	366	C26H54	013475-76-8	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP081021\
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Sample : M3208-06
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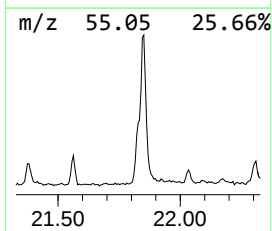
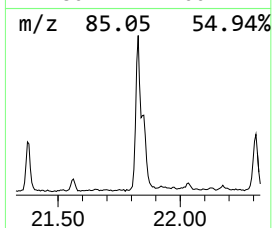
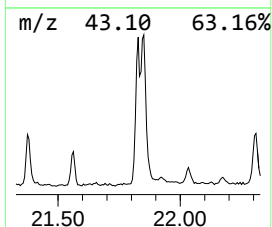
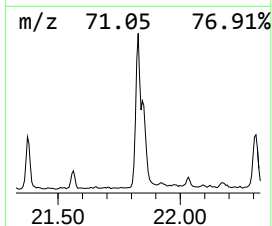
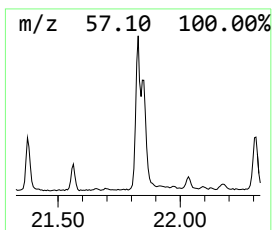
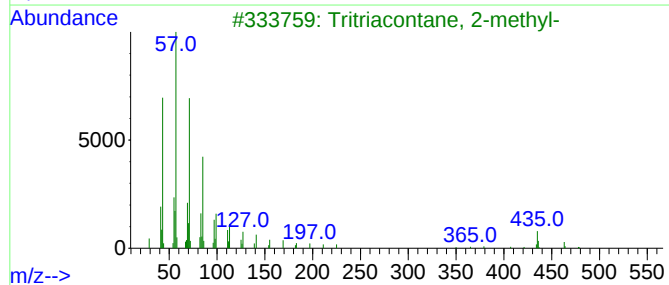
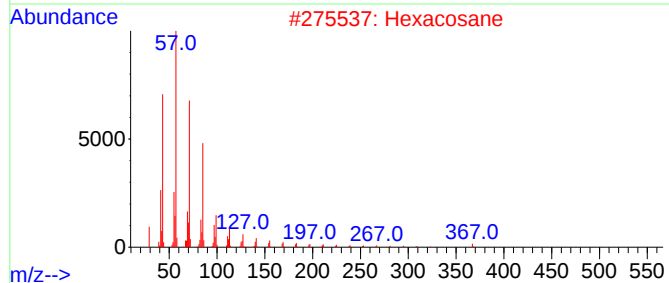
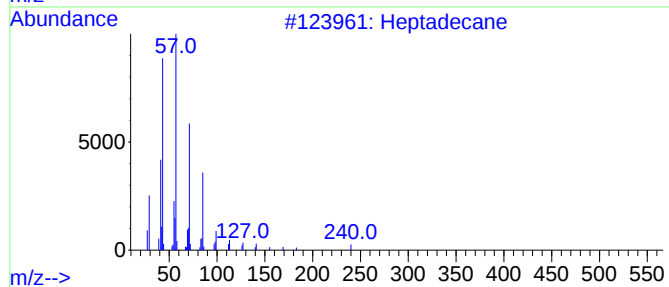
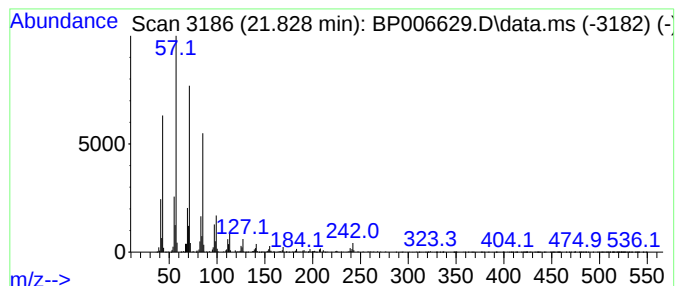
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 (DEL) Alkane: Straight-Chai... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD		R.T.	
21.830	5.78 ng/ul	724392	Chrysene-d12		21.216	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Heptadecane		240	C17H36	000629-78-7	93
2	Hexacosane		366	C26H54	000630-01-3	91
3	Tritriacontane, 2-methyl-		479	C34H70	066214-27-5	91
4	Octacosane		394	C28H58	000630-02-4	91
5	Hexatriacontane		507	C36H74	000630-06-8	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP081021\
Data File : BP006629.D
Acq On : 10 Aug 2021 18:25
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Sample : M3208-06
Misc :
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ClientSampleId :
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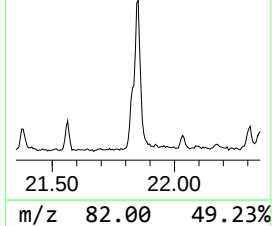
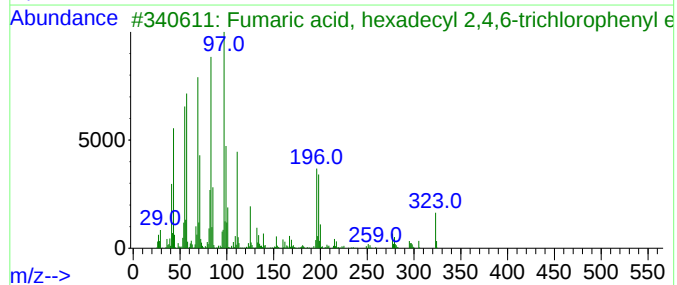
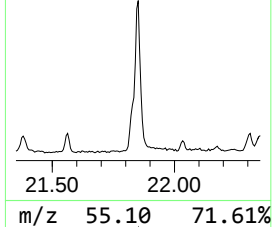
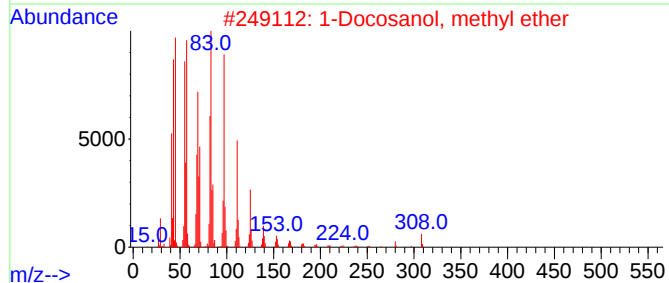
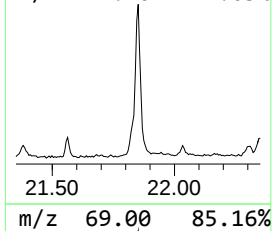
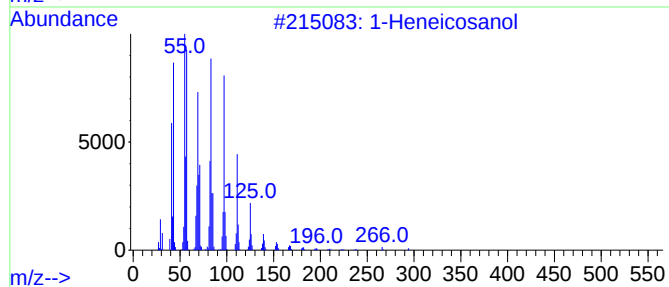
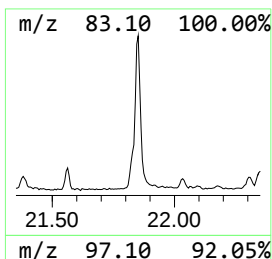
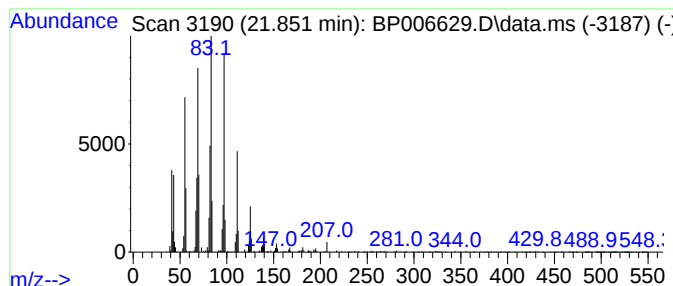
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 1-Heneicosanol Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.850	11.50 ng/ul	1441120	Chrysene-d12	21.216

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Heneicosanol	312	C21H44O	015594-90-8	90
2			1-Docosanol, methyl ether	340	C23H48O	1000333-91-9	81
3			Fumaric acid, hexadecyl 2,4,6-tr...	518	C26H37Cl3O4	1000348-27-9	80
4			Fumaric acid, 2-chloroethyl hexa...	402	C22H39ClO4	1000330-62-2	72
5			Fumaric acid, 3-fluorophenyl hex...	434	C26H39FO4	1000345-21-2	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP081021\
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Acq On : 10 Aug 2021 18:25
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Sample : M3208-06
Misc :
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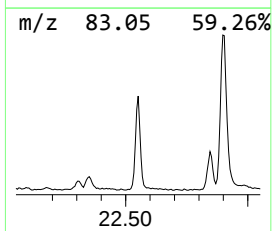
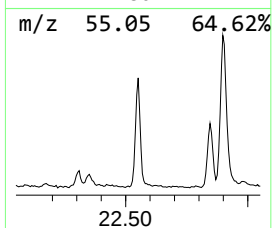
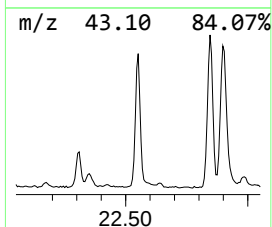
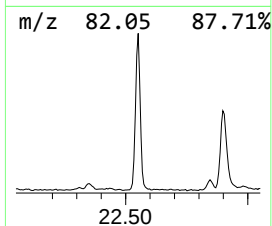
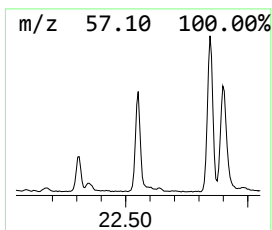
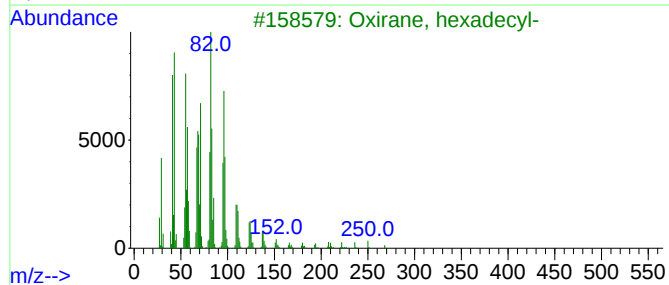
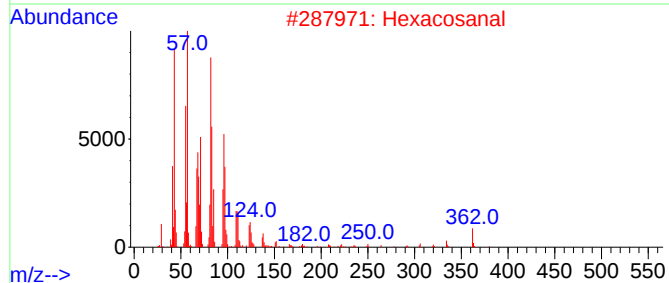
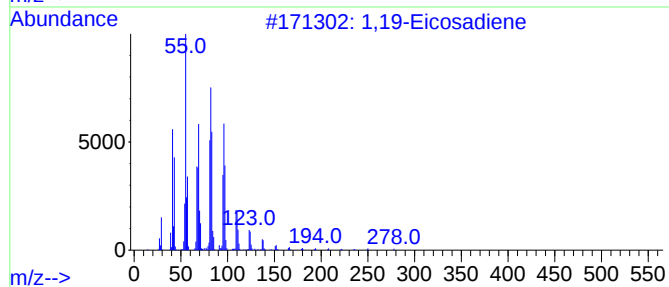
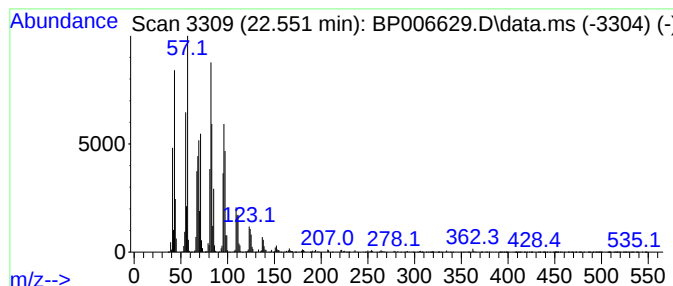
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 1,19-Eicosadiene Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.550	14.06 ng/ul	1794110	Perylene-d12	23.527

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,19-Eicosadiene			278	C20H38	014811-95-1	95
2	Hexacosanal			380	C26H52O	026627-85-0	94
3	Oxirane, hexadecyl-			268	C18H36O	007390-81-0	94
4	Oxirane, heptadecyl-			282	C19H38O	067860-04-2	91
5	Oxirane, hexadecyl-			268	C18H36O	007390-81-0	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP081021\
Data File : BP006629.D
Acq On : 10 Aug 2021 18:25
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Sample : M3208-06
Misc :
ALS Vial : 14 Sample Multiplier: 1

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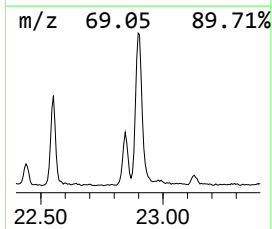
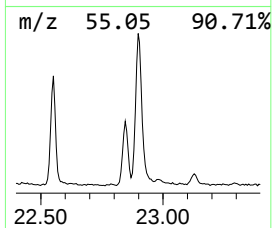
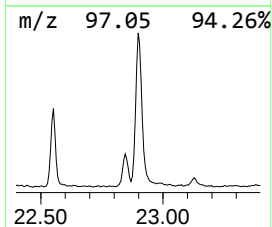
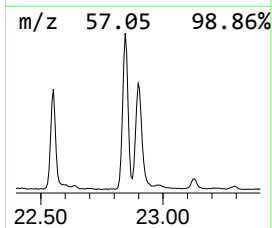
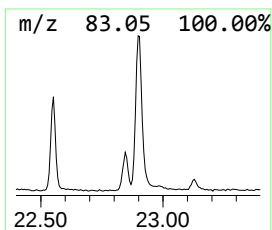
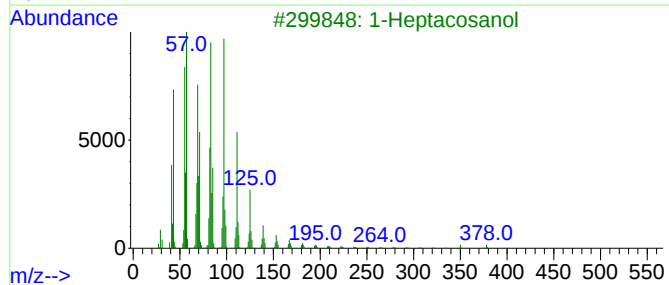
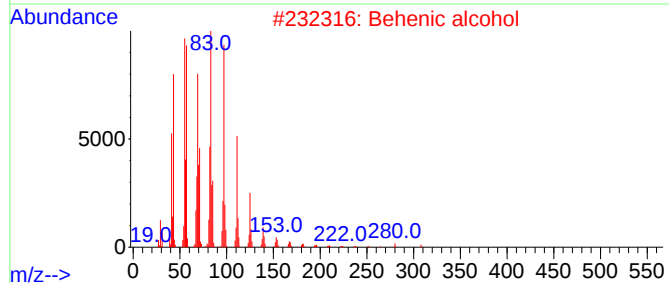
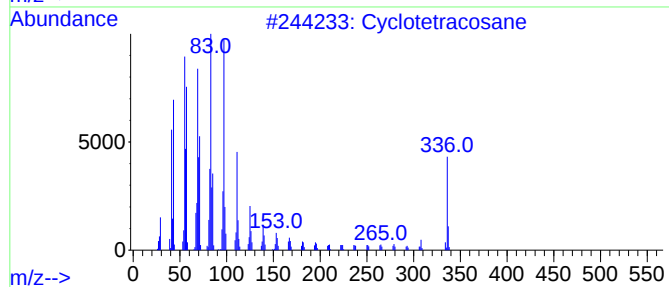
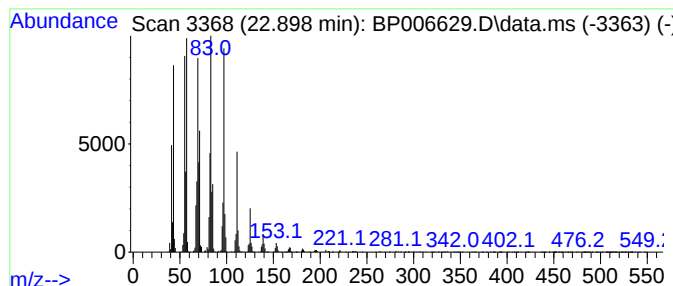
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TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 8 (DEL) Alkane: Cyclic22.90 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.900	19.49 ng/ul	2487060	Perylene-d12	23.527

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclotetracosane	336	C24H48	000297-03-0	99
2			Behenic alcohol	326	C22H46O	000661-19-8	94
3			1-Heptacosanol	396	C27H56O	002004-39-9	94
4			1-Heneicosanol	312	C21H44O	015594-90-8	93
5			Octacosanol	410	C28H58O	000557-61-9	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP081021\
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ClientSampleId :
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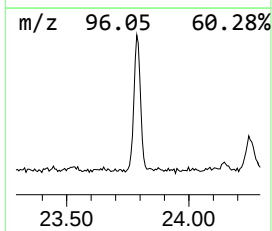
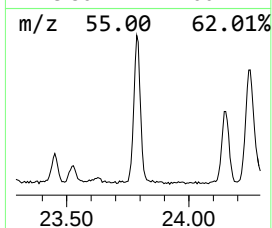
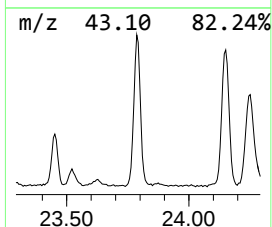
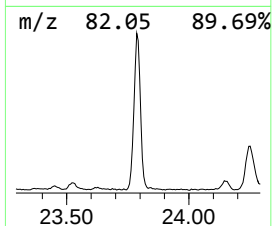
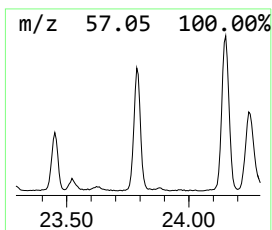
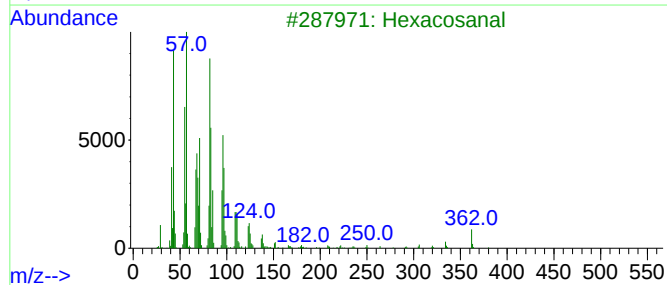
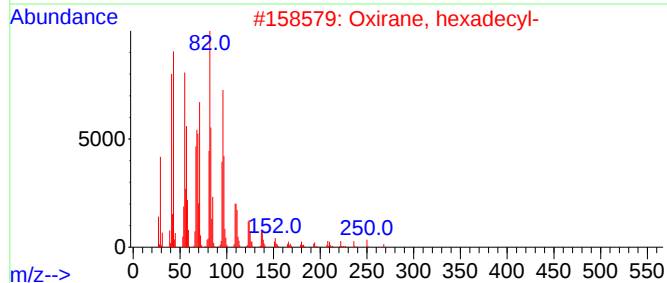
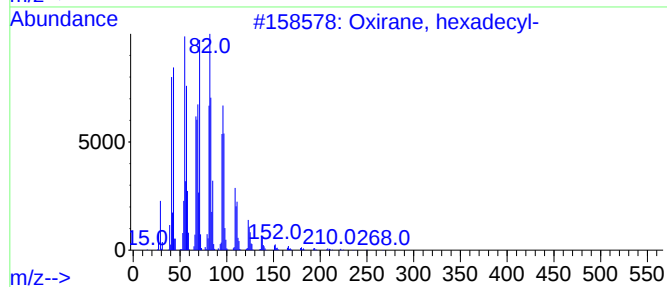
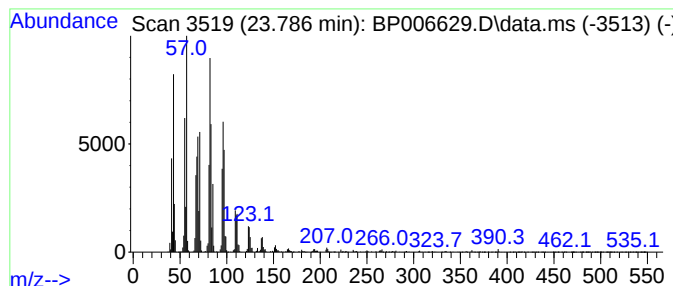
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 Oxirane, hexadecyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.790	17.36 ng/ul	2214480	Perylene-d12	23.527

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Oxirane, hexadecyl-	268	C18H36O	007390-81-0	94		
2	Oxirane, hexadecyl-	268	C18H36O	007390-81-0	94		
3	Hexacosanal	380	C26H52O	026627-85-0	94		
4	Oxirane, heptadecyl-	282	C19H38O	067860-04-2	94		
5	Hexadecanal	240	C16H32O	000629-80-1	93		



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP081021\
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Acq On : 10 Aug 2021 18:25
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Misc :
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ClientSampleId :
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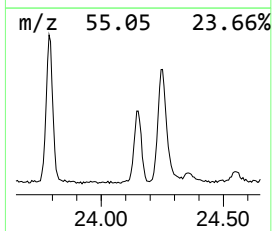
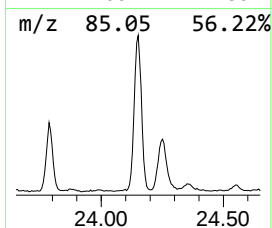
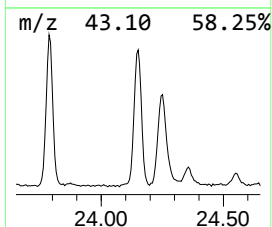
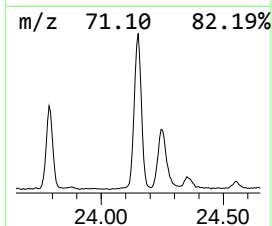
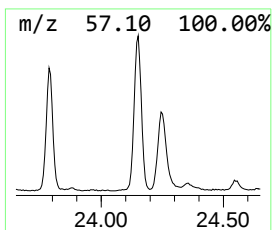
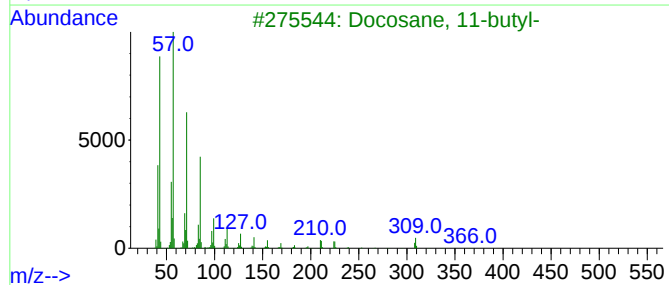
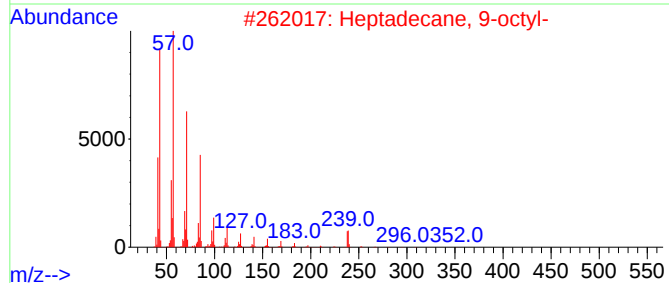
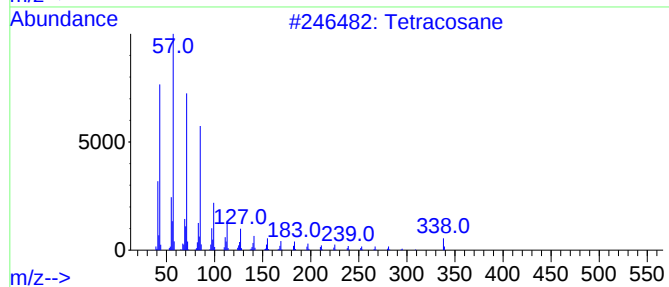
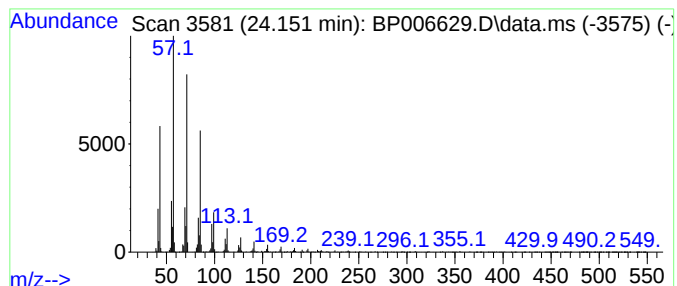
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 (DEL) Alkane: Straight-Chai... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.150	11.73 ng/ul	1497000	Perylene-d12	23.527

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tetracosane	338	C24H50	000646-31-1	97
2			Heptadecane, 9-octyl-	352	C25H52	007225-64-1	94
3			Docosane, 11-butyl-	366	C26H54	013475-76-8	94
4			2-Methylpentacosane	366	C26H54	000629-87-8	94
5			Docosane, 9-butyl-	366	C26H54	055282-14-9	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP081021\
Data File : BP006629.D
Acq On : 10 Aug 2021 18:25
Operator : CG/JU
Sample : M3208-06
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C00E6

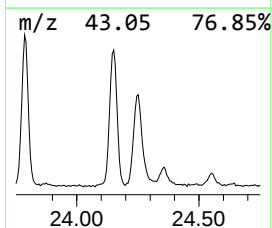
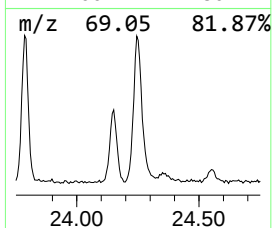
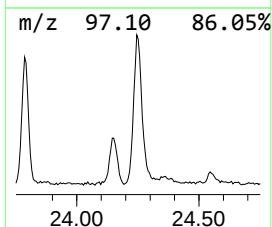
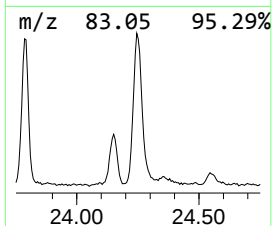
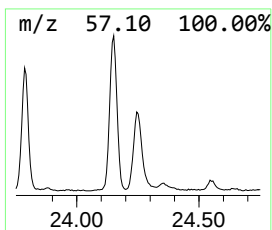
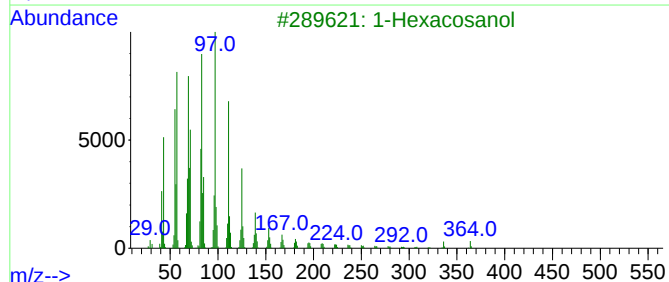
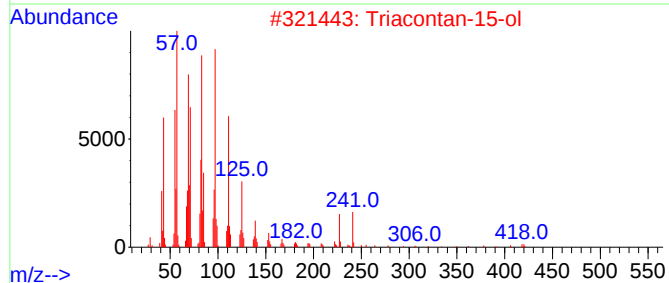
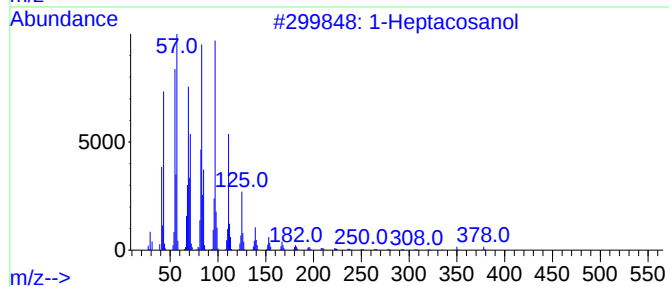
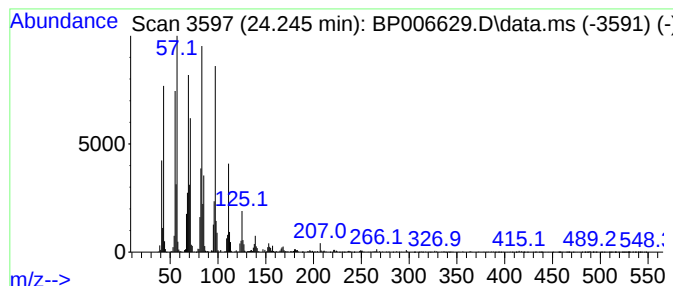
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Quant Title : SVOA CALIBRATION

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

Peak Number 11 1-Heptacosanol Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.250	13.72 ng/ul	1750510	Perylene-d12	23.527

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Heptacosanol	396	C27H56O	002004-39-9	93
2			Triacontan-15-ol	438	C30H62O	031849-26-0	91
3			1-Hexacosanol	382	C26H54O	000506-52-5	91
4			Octacosanol	410	C28H58O	000557-61-9	91
5			Heptacosyl acetate	438	C29H58O2	1000351-78-2	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP081021\
 Data File : BP006629.D
 Acq On : 10 Aug 2021 18:25
 Operator : CG/JU
 Sample : M3208-06
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C00E6

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP081021.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-(2-b...	10.300	10.9	ng/ul	813066	2	10.505	1493710	20.0
n-Hexadecanoic ...	18.020	5.0	ng/ul	540623	4	17.116	2139640	20.0
(DEL) Alkane: C...	20.950	13.6	ng/ul	1699960	5	21.216	2505530	20.0
(DEL) Alkane: S...	21.370	2.0	ng/ul	254316	5	21.216	2505530	20.0
(DEL) Alkane: S...	21.830	5.8	ng/ul	724392	5	21.216	2505530	20.0
1-Heneicosanol	21.850	11.5	ng/ul	1441120	5	21.216	2505530	20.0
1,19-Eicosadiene	22.550	14.1	ng/ul	1794110	6	23.527	2551700	20.0
(DEL) Alkane: C...	22.900	19.5	ng/ul	2487060	6	23.527	2551700	20.0
Oxirane, hexade...	23.790	17.4	ng/ul	2214480	6	23.527	2551700	20.0
(DEL) Alkane: S...	24.150	11.7	ng/ul	1497000	6	23.527	2551700	20.0
1-Heptacosanol	24.250	13.7	ng/ul	1750510	6	23.527	2551700	20.0