

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP050824\
 Data File : BP020264.D
 Acq On : 09 May 2024 10:45
 Operator : MA/JU
 Sample : P2369-08
 Misc :
 ALS Vial : 32 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 A43N2

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-BP050124.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BP020264.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.217	19	22	33	rVB	22850	38532	0.75%	0.075%
2	3.481	63	67	82	rBV	86170	153135	3.00%	0.298%
3	3.617	86	90	105	rBV	271356	477804	9.35%	0.931%
4	3.911	137	140	144	rVB2	6809	7526	0.15%	0.015%
5	6.005	492	496	506	rBV5	4799	10069	0.20%	0.020%
6	6.664	603	608	614	rBV9	4667	10876	0.21%	0.021%
7	6.805	624	632	654	rBV	363413	723585	14.17%	1.410%
8	6.969	654	660	673	rVV	330954	558589	10.94%	1.088%
9	7.158	685	692	701	rVV	411859	738468	14.46%	1.439%
10	7.258	704	709	727	rVB	41501	106700	2.09%	0.208%
11	7.616	763	770	784	rBV	410720	688298	13.48%	1.341%
12	7.716	784	787	792	rVB5	5246	8379	0.16%	0.016%
13	8.328	883	891	911	rBV2	434054	862980	16.89%	1.681%
14	8.775	960	967	980	rBV	466641	826309	16.18%	1.610%
15	8.887	982	986	990	rVV6	6382	13788	0.27%	0.027%
16	8.934	990	994	1000	rVB9	5642	10390	0.20%	0.020%
17	9.352	1060	1065	1074	rBV	33000	62639	1.23%	0.122%
18	9.487	1081	1088	1105	rBV	383110	747250	14.63%	1.456%
19	9.758	1129	1134	1135	rBV5	5883	6906	0.14%	0.013%
20	10.022	1172	1179	1202	rBV	535314	1159026	22.69%	2.258%
21	10.387	1232	1241	1255	rBV	625754	1048051	20.52%	2.042%
22	10.552	1261	1269	1291	rBV	476082	1056389	20.68%	2.058%
23	10.975	1335	1341	1353	rBV	9264	24991	0.49%	0.049%
24	11.163	1364	1373	1381	rBV2	64363	168807	3.30%	0.329%
25	11.940	1500	1505	1509	rBV6	3728	7625	0.15%	0.015%
26	11.993	1509	1514	1521	rVB2	9467	18532	0.36%	0.036%
27	12.581	1608	1614	1617	rBV7	6270	12164	0.24%	0.024%
28	12.875	1660	1664	1668	rBV7	3494	5883	0.12%	0.011%
29	13.081	1693	1699	1703	rBV8	2884	6269	0.12%	0.012%
30	13.169	1707	1714	1722	rVV	276606	508140	9.95%	0.990%
31	13.246	1723	1727	1733	rVB8	14985	24642	0.48%	0.048%
32	13.704	1794	1805	1819	rBV	1376748	1934827	37.88%	3.769%
33	13.981	1841	1852	1866	rBV2	1373210	2239094	43.84%	4.362%
34	14.293	1898	1905	1914	rBV	993028	1515376	29.67%	2.952%
35	14.510	1933	1942	1943	rBV	2205391	2359800	46.20%	4.597%
36	14.534	1943	1946	1974	rVV3	2770613	5107911	100.00%	9.951%

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Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP050124.MA.M
Title : SVOA CALIBRATION

37	14.745	1974	1982	1988	rVB5	29817	64476	1.26%	0.126%
38	14.951	2013	2017	2023	rVB2	23924	36331	0.71%	0.071%
39	15.116	2040	2045	2049	rVB3	5093	7743	0.15%	0.015%
40	15.169	2049	2054	2061	rVB7	6158	12238	0.24%	0.024%
41	15.257	2064	2069	2072	rBV5	6572	10877	0.21%	0.021%
42	15.316	2072	2079	2087	rBV	1868138	2822333	55.25%	5.498%
43	15.475	2099	2106	2117	rBV	383718	676290	13.24%	1.317%
44	15.675	2135	2140	2143	rBV6	8189	12618	0.25%	0.025%
45	15.904	2175	2179	2186	rBV7	18653	35564	0.70%	0.069%
46	15.957	2186	2188	2194	rVB6	3694	5628	0.11%	0.011%
47	16.075	2202	2208	2209	rBV6	7087	10226	0.20%	0.020%
48	16.228	2231	2234	2236	rBV4	6030	8444	0.17%	0.016%
49	16.251	2236	2238	2242	rVB5	8778	10594	0.21%	0.021%
50	16.316	2245	2249	2255	rVB	18458	29820	0.58%	0.058%
51	16.404	2258	2264	2265	rBV6	6144	10236	0.20%	0.020%
52	16.534	2281	2286	2293	rBV2	69497	111979	2.19%	0.218%
53	16.669	2305	2309	2315	rVB9	10300	18329	0.36%	0.036%
54	16.769	2319	2326	2337	rBV	2095263	3258202	63.79%	6.347%
55	16.922	2348	2352	2368	rVB10	33928	108815	2.13%	0.212%
56	17.069	2372	2377	2381	rVV5	23794	35395	0.69%	0.069%
57	17.134	2381	2388	2393	rVV	1118918	1911636	37.43%	3.724%
58	17.175	2393	2395	2400	rVV2	130026	204559	4.00%	0.398%
59	17.234	2400	2405	2417	rVB2	1955663	2956702	57.88%	5.760%
60	17.328	2417	2421	2426	rBV2	38616	54082	1.06%	0.105%
61	17.428	2435	2438	2445	rVB8	8144	18100	0.35%	0.035%
62	17.563	2455	2461	2464	rBV4	21351	42087	0.82%	0.082%
63	17.675	2475	2480	2485	rBV7	10266	27108	0.53%	0.053%
64	17.769	2492	2496	2500	rVB2	57290	81232	1.59%	0.158%
65	17.857	2505	2511	2516	rVB3	39985	68180	1.33%	0.133%
66	17.922	2519	2522	2525	rVV	24194	34248	0.67%	0.067%
67	17.957	2525	2528	2537	rVB4	32896	56770	1.11%	0.111%
68	18.092	2544	2551	2565	rBV	1670082	2576850	50.45%	5.020%
69	18.245	2573	2577	2582	rBV2	21984	35872	0.70%	0.070%
70	18.392	2599	2602	2605	rBV2	23796	27863	0.55%	0.054%
71	18.434	2605	2609	2620	rVB8	31967	92038	1.80%	0.179%
72	18.586	2630	2635	2641	rBV5	32700	62850	1.23%	0.122%
73	18.675	2646	2650	2657	rVB	79653	111268	2.18%	0.217%
74	18.786	2666	2669	2674	rBV7	14266	23267	0.46%	0.045%
75	18.922	2690	2692	2697	rVB6	11584	10709	0.21%	0.021%
76	19.010	2702	2707	2712	rBV4	36165	75205	1.47%	0.147%

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Title : SVOA CALIBRATION

77	19.063	2712	2716	2717	rVV3	26624	33800	0.66%	0.066%
78	19.086	2717	2720	2727	rVB4	87490	123918	2.43%	0.241%
79	19.175	2732	2735	2741	rVB4	40241	51104	1.00%	0.100%
80	19.228	2741	2744	2746	rBV	32165	36903	0.72%	0.072%
81	19.281	2749	2753	2757	rBV	239213	339009	6.64%	0.660%
82	19.439	2775	2780	2791	rBV2	786439	1258026	24.63%	2.451%
83	19.586	2801	2805	2808	rBV	87725	109282	2.14%	0.213%
84	19.633	2808	2813	2825	rVB	2516263	3702894	72.49%	7.214%
85	19.928	2859	2863	2869	rBV5	65835	118356	2.32%	0.231%
86	21.310	3094	3098	3107	rVB	311118	537289	10.52%	1.047%
87	21.633	3146	3153	3159	rBV	995664	1803354	35.31%	3.513%
88	24.763	3676	3685	3696	rBV2	923752	2635668	51.60%	5.134%
89	25.004	3717	3726	3737	rVB	555249	1576669	30.87%	3.071%

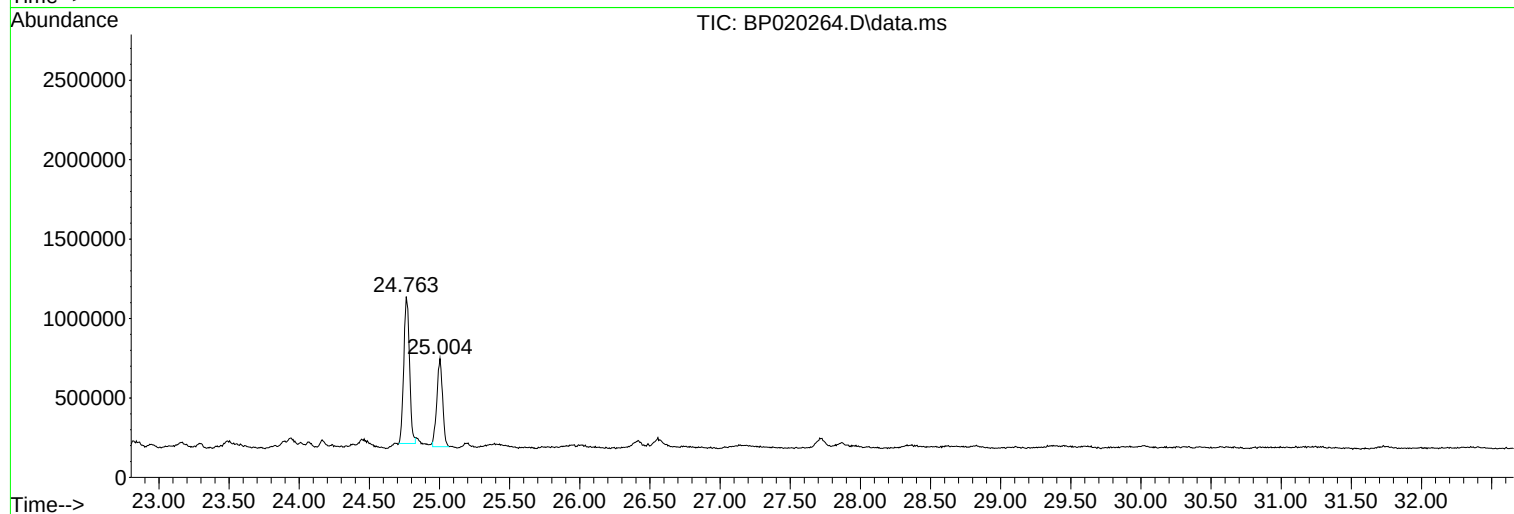
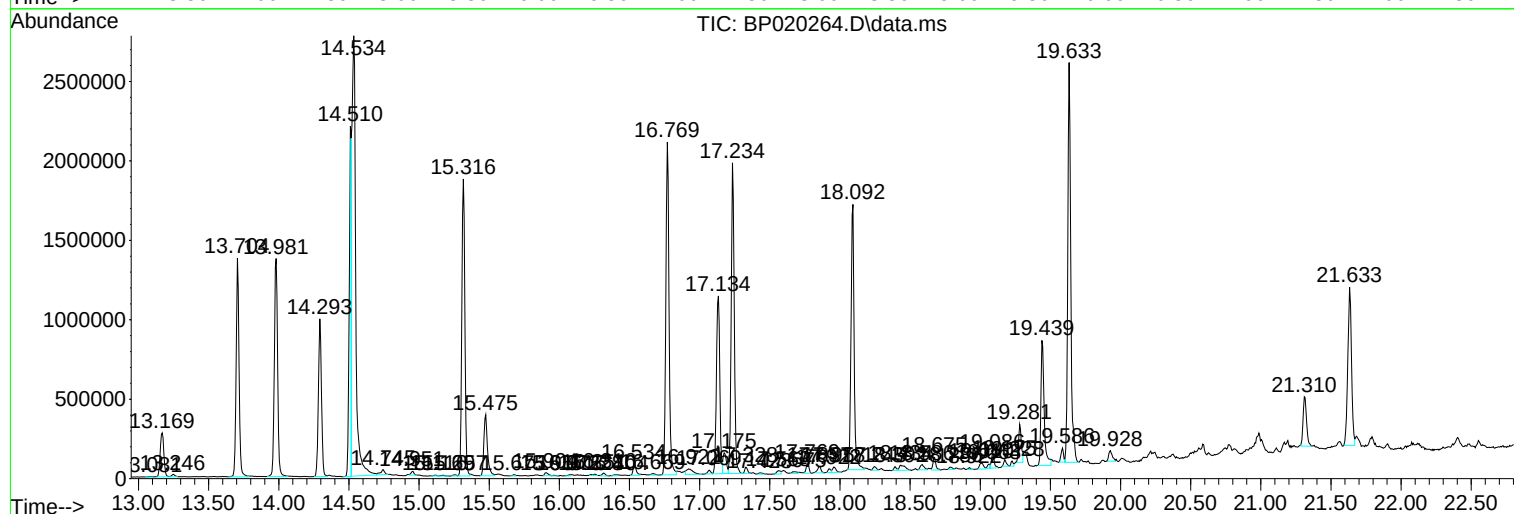
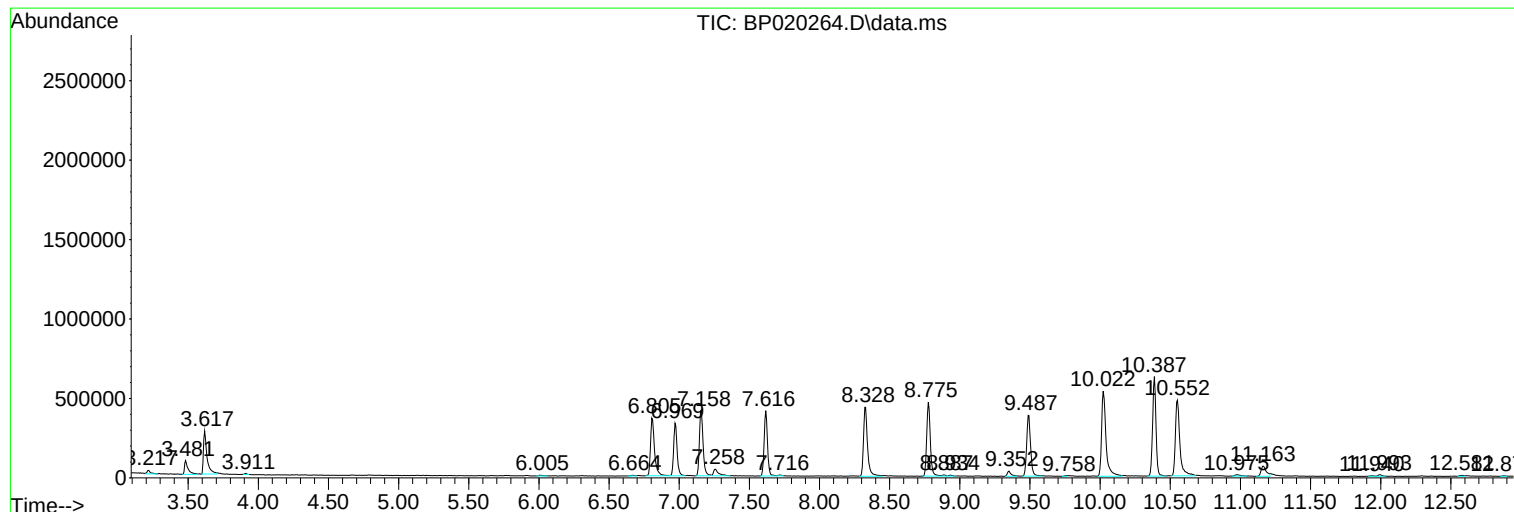
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP050124.MA.M
Quant Title : SVOA CALIBRATION

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



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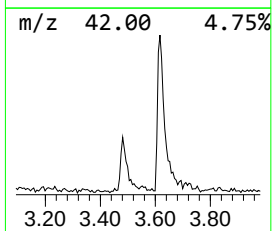
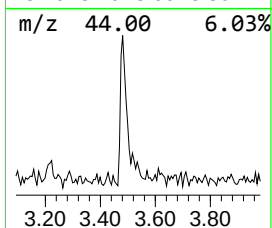
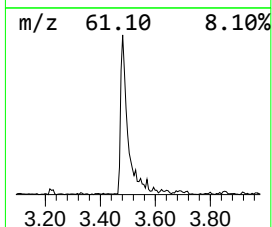
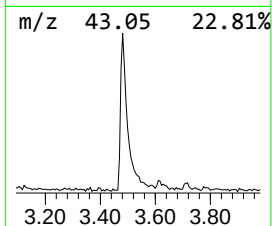
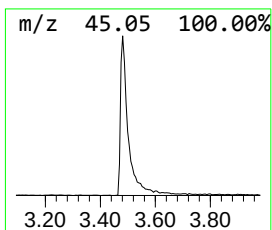
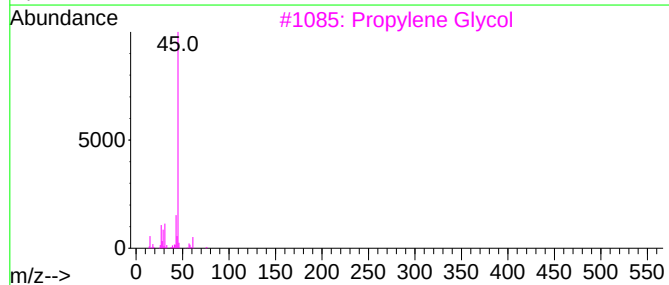
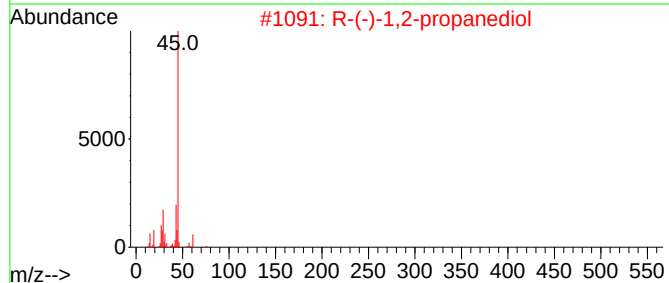
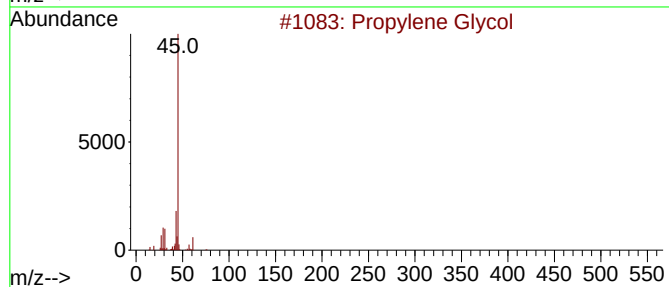
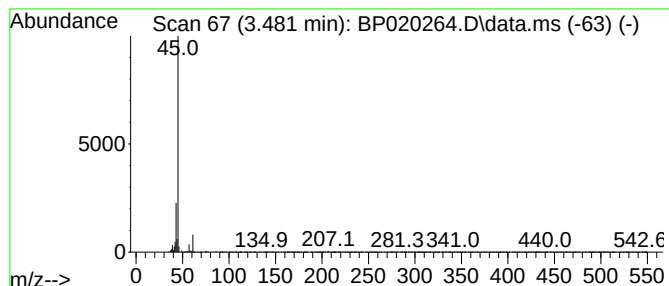
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP050124.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Propylene Glycol Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.481	4.45 ng/ul	153135	1,4-Dichlorobenzene-d4	7.616

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Propylene Glycol	76	C3H8O2	000057-55-6	80
2			R-(-)-1,2-propanediol	76	C3H8O2	004254-14-2	56
3			Propylene Glycol	76	C3H8O2	000057-55-6	56
4			Isopropyl Alcohol	60	C3H8O	000067-63-0	50
5			Isopropyl Alcohol	60	C3H8O	000067-63-0	9



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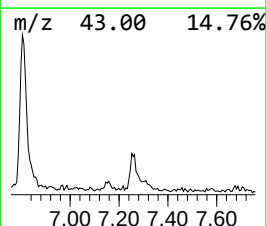
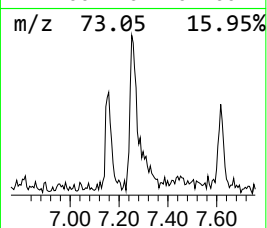
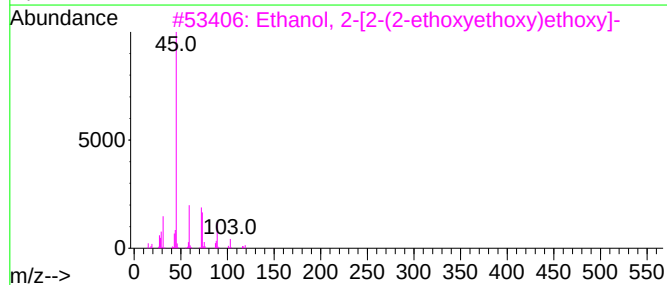
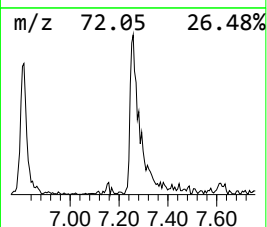
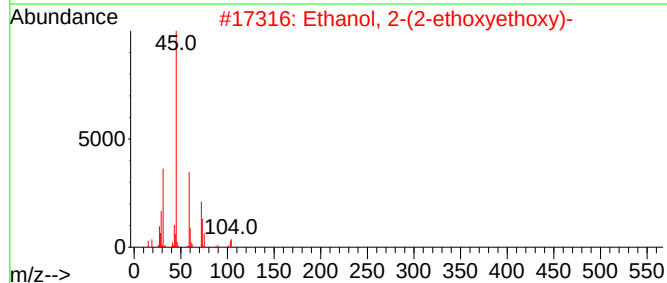
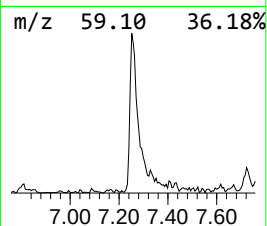
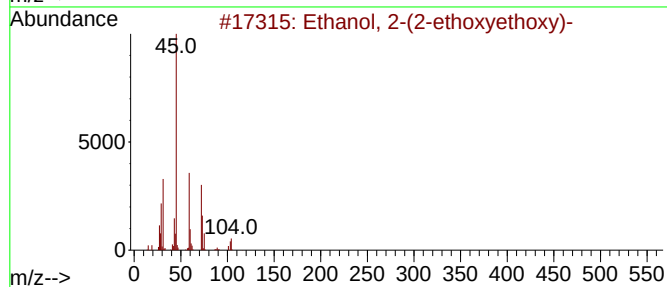
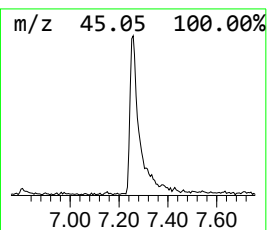
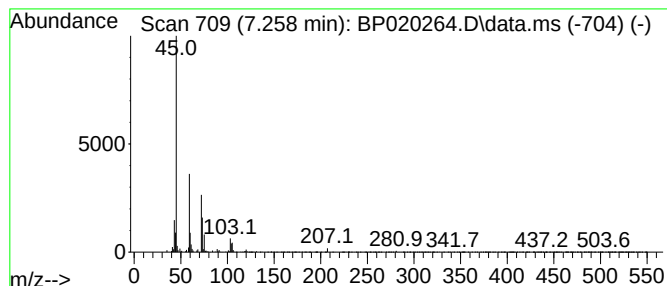
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP050124.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 Ethanol, 2-(2-ethoxyethoxy)- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.258	3.10 ng/ul	106700	1,4-Dichlorobenzene-d4	7.616

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Ethanol, 2-(2-ethoxyethoxy)-	134	C6H14O3	000111-90-0	90
2			Ethanol, 2-(2-ethoxyethoxy)-	134	C6H14O3	000111-90-0	83
3			Ethanol, 2-[2-(2-ethoxyethoxy)et...	178	C8H18O4	000112-50-5	78
4			Ethanol, 2-(2-ethoxyethoxy)-	134	C6H14O3	000111-90-0	78
5			Diethyl carbitol	162	C8H18O3	000112-36-7	72



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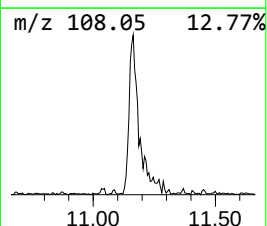
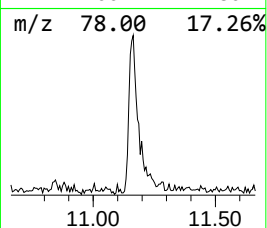
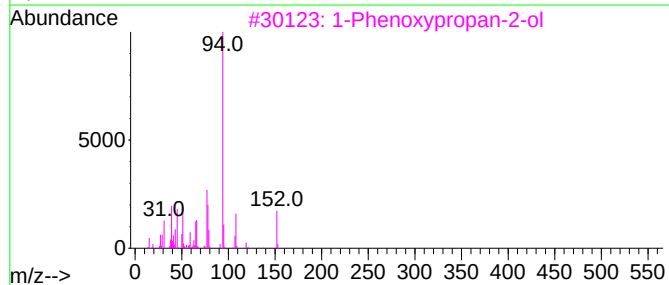
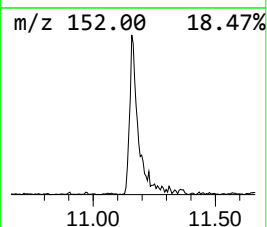
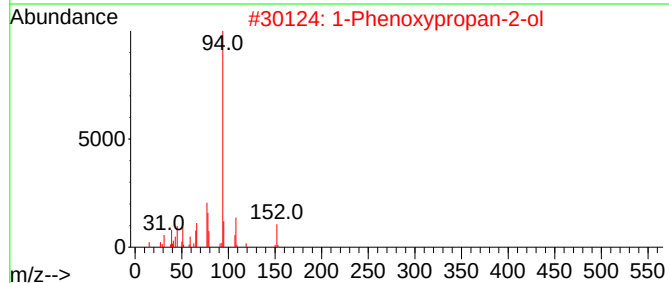
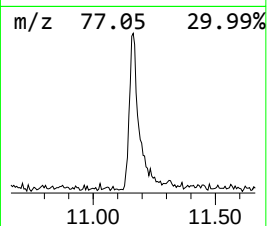
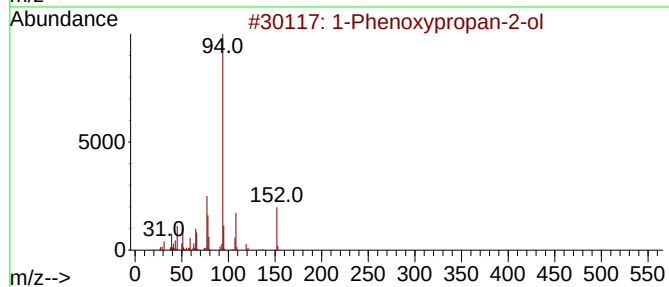
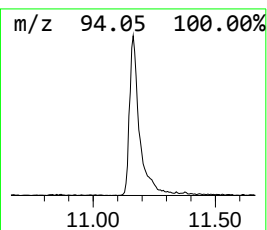
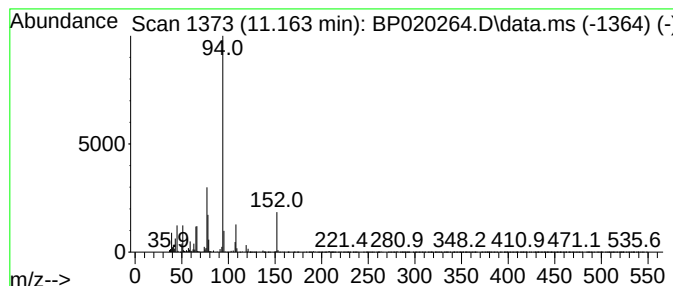
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP050124.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 1-Phenoxypropan-2-ol Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.163	3.22 ng/ul	168807	Naphthalene-d8	10.387

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	95
2			1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	93
3			1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	91
4			1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	87
5			1-Propanol, 3-phenoxy-	152	C9H12O2	006180-61-6	64



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ClientSampleId :
A43N2

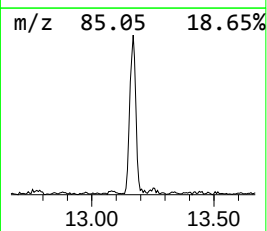
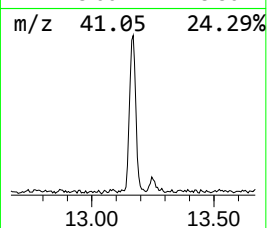
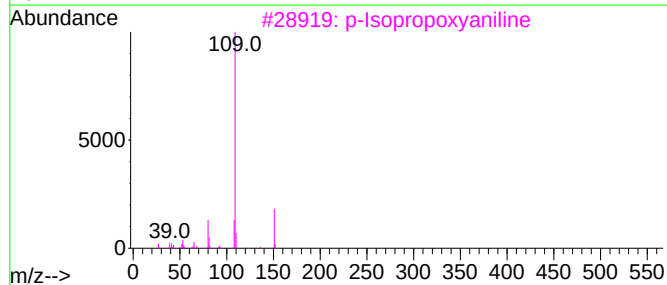
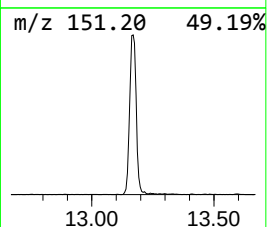
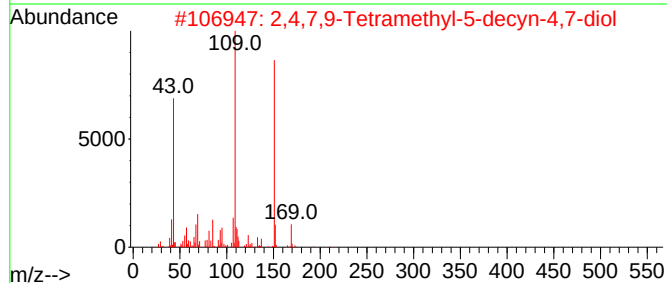
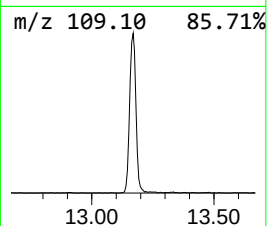
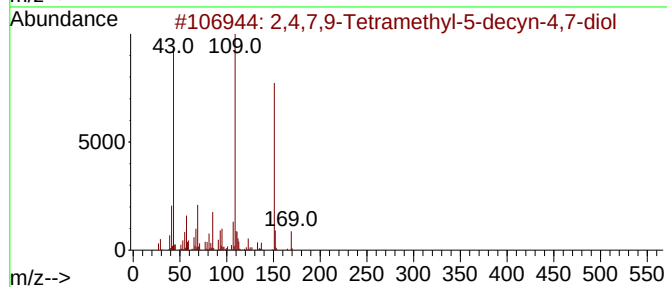
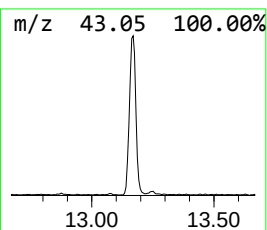
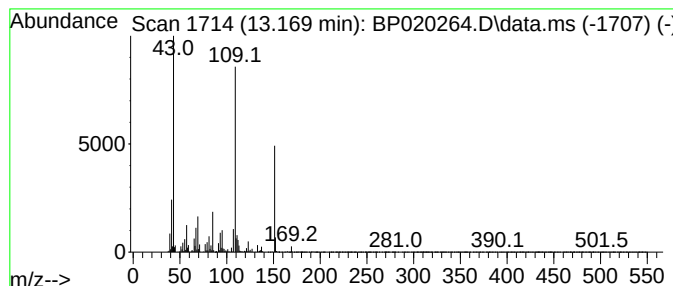
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 2,4,7,9-Tetramethyl-5-decyn... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.169	6.71 ng/ul	508140	Acenaphthene-d10	14.293

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,4,7,9-Tetramethyl-5-decyn-4,7-...	226	C14H26O2	000126-86-3	91		
2	2,4,7,9-Tetramethyl-5-decyn-4,7-...	226	C14H26O2	000126-86-3	90		
3	p-Isopropoxyaniline	151	C9H13NO	007664-66-6	64		
4	Pyridine, 2-butyl-, 1-oxide	151	C9H13NO	031396-32-4	59		
5	Acetaminophen	151	C8H9NO2	000103-90-2	59		



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP050824\
Data File : BP020264.D
Acq On : 09 May 2024 10:45
Operator : MA/JU
Sample : P2369-08
Misc :
ALS Vial : 32 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A43N2

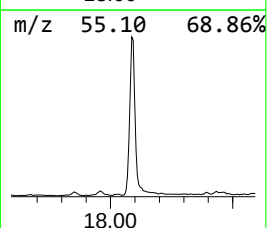
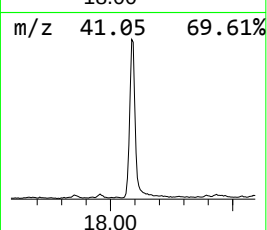
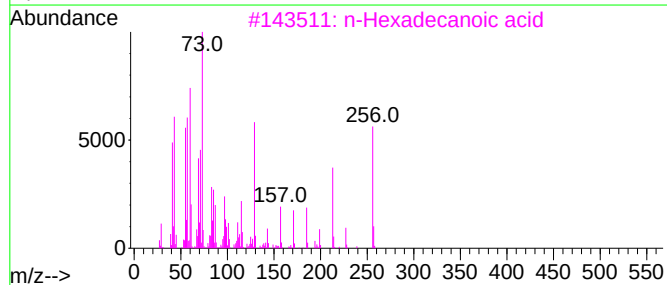
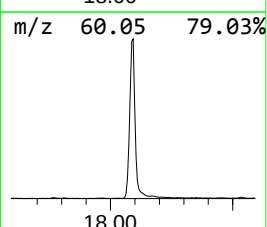
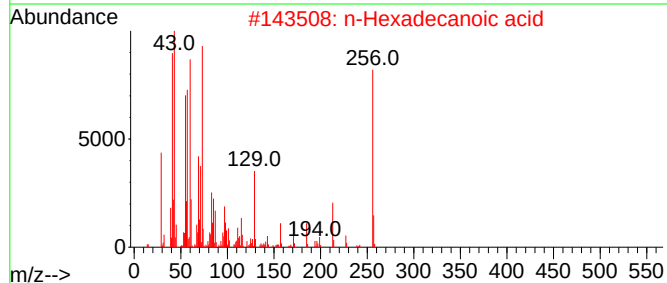
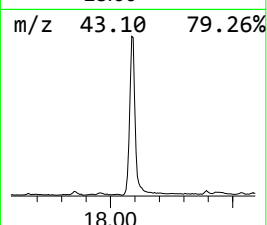
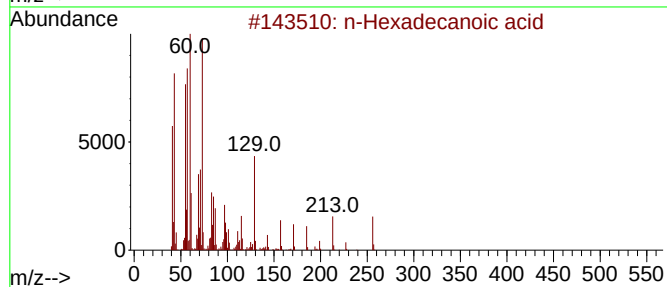
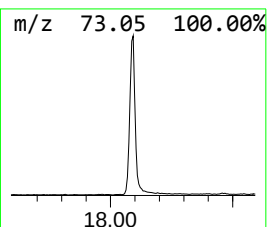
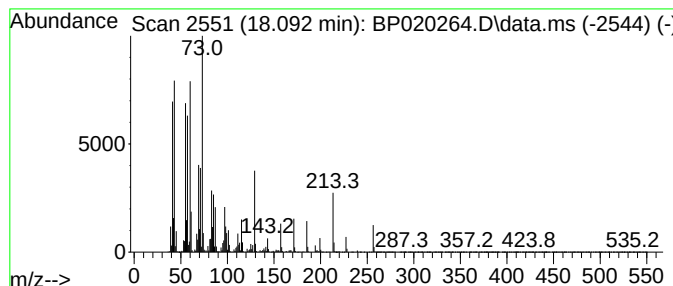
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 n-Hexadecanoic acid Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.092	26.96 ng/ul	2576850	Phenanthrene-d10	17.134

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	99
3			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
4			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
5			Tetradecanoic acid	228	C14H28O2	000544-63-8	95



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP050824\
Data File : BP020264.D
Acq On : 09 May 2024 10:45
Operator : MA/JU
Sample : P2369-08
Misc :
ALS Vial : 32 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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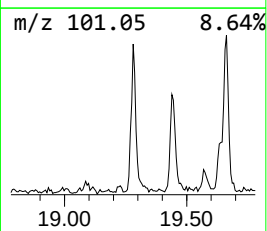
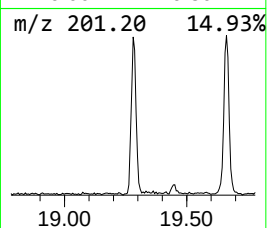
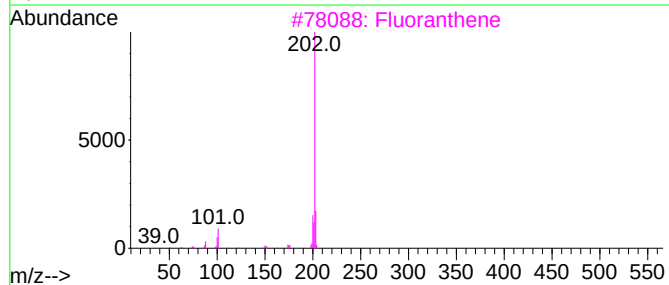
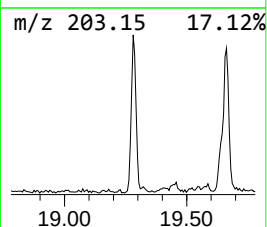
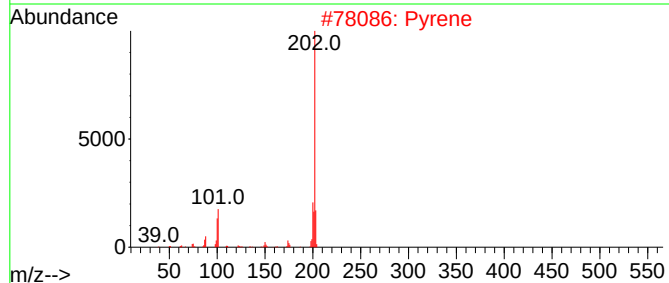
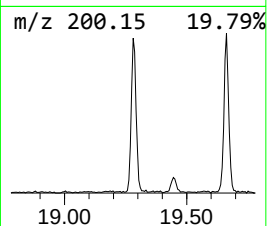
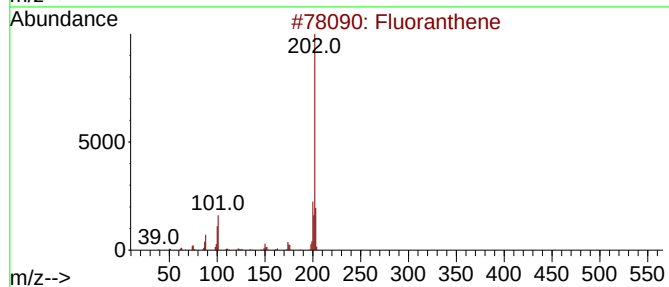
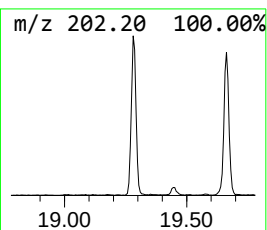
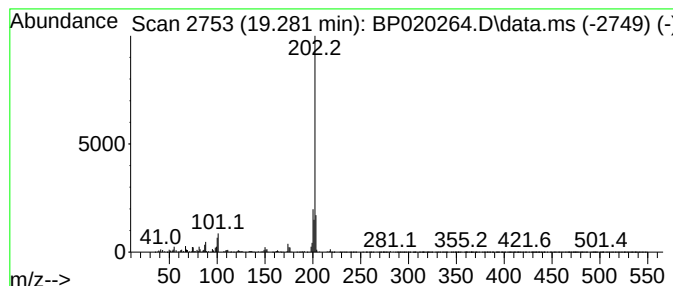
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 Fluoranthene Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.281	3.55 ng/ul	339009	Phenanthrene-d10	17.134

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Fluoranthene	202	C16H10	000206-44-0	95
2			Pyrene	202	C16H10	000129-00-0	90
3			Fluoranthene	202	C16H10	000206-44-0	87
4			Fluoranthene	202	C16H10	000206-44-0	83
5			Pyrene	202	C16H10	000129-00-0	81



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP050824\
Data File : BP020264.D
Acq On : 09 May 2024 10:45
Operator : MA/JU
Sample : P2369-08
Misc :
ALS Vial : 32 Sample Multiplier: 1

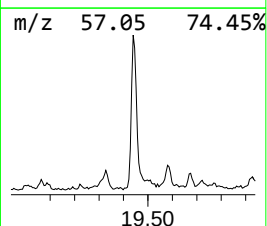
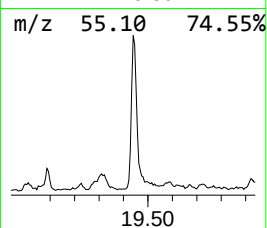
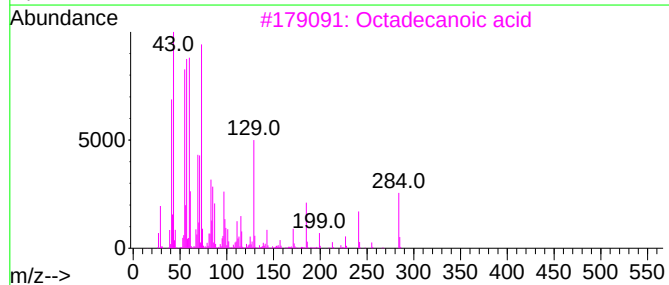
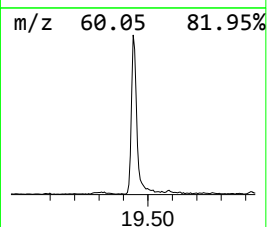
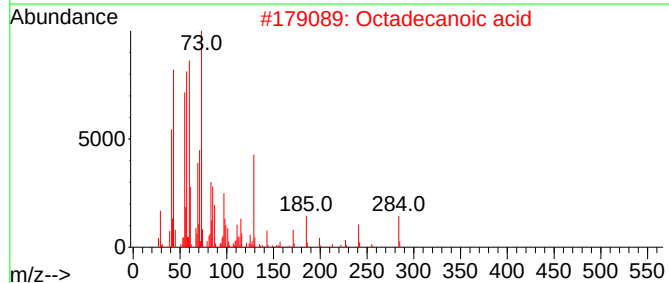
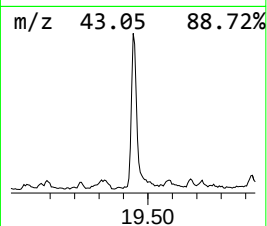
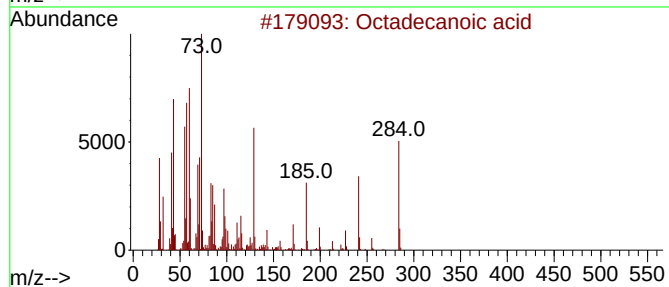
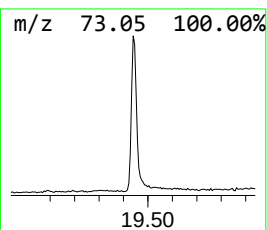
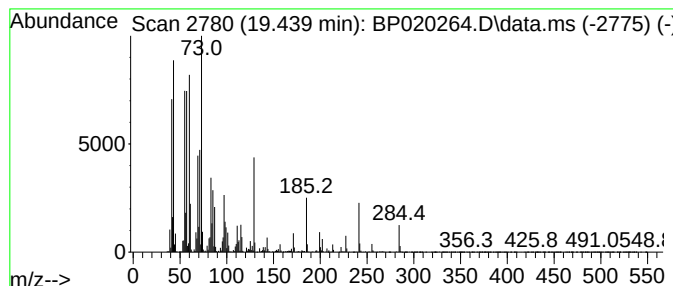
Instrument :
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP050124.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 7 Octadecanoic acid Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD		R.T.	
19.439	13.95 ng/ul	1258030	Chrysene-d12		21.633	
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	96
5	Octadecanoic acid		284	C18H36O2	000057-11-4	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP050824\
Data File : BP020264.D
Acq On : 09 May 2024 10:45
Operator : MA/JU
Sample : P2369-08
Misc :
ALS Vial : 32 Sample Multiplier: 1

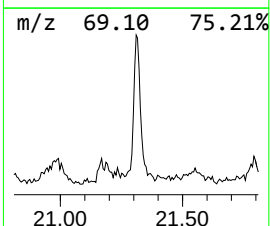
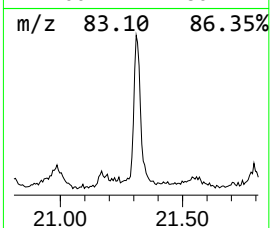
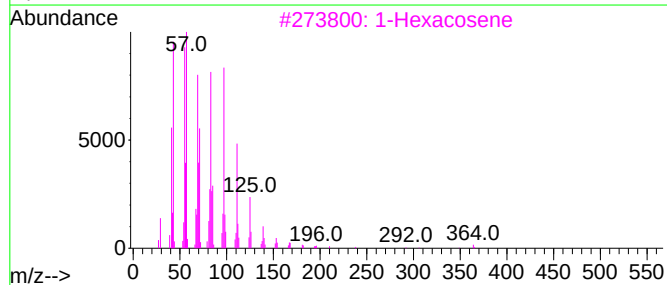
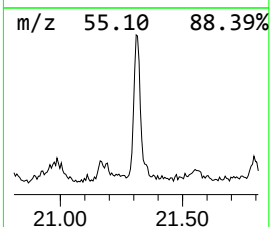
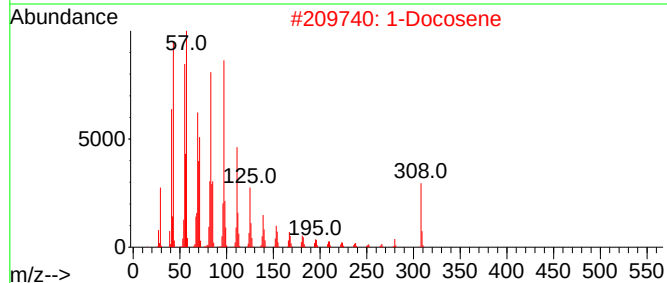
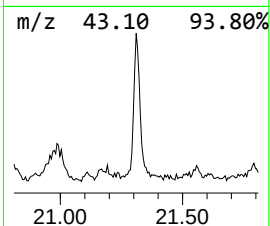
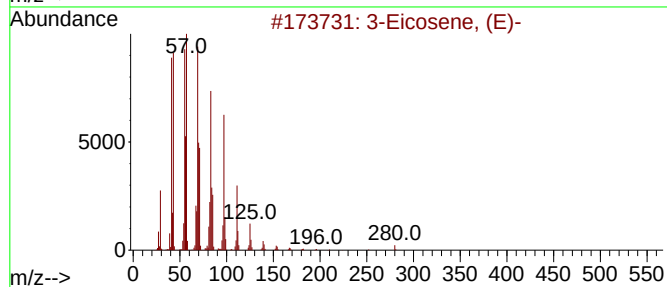
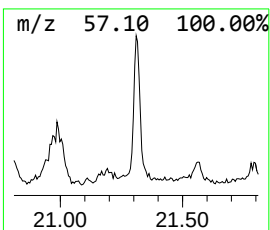
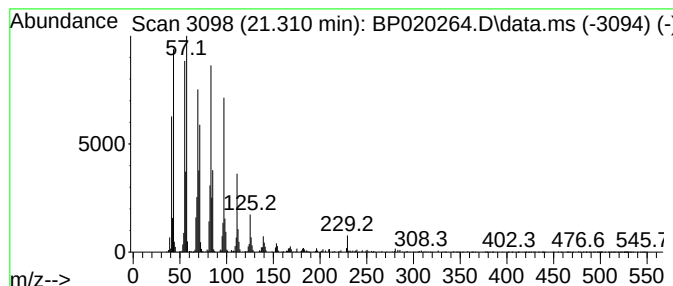
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP050124.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 8 (DEL) Alkane: Straight-Chai... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD		R.T.	
21.310	5.96 ng/ul	537289	Chrysene-d12		21.633	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	3-Eicosene, (E)-		280	C20H40	074685-33-9	97
2	1-Docosene		308	C22H44	001599-67-3	95
3	1-Hexacosene		364	C26H52	018835-33-1	95
4	5-Octadecene, (E)-		252	C18H36	007206-21-5	94
5	Heptadecyl heptafluorobutyrate		452	C21H35F7O2	959085-66-6	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP050824\
Data File : BP020264.D
Acq On : 09 May 2024 10:45
Operator : MA/JU
Sample : P2369-08
Misc :
ALS Vial : 32 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP050124.MA.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
Propylene Glycol	3.481	4.5	ng/ul	153135	1	7.616	688298	20.0
Ethanol, 2-(2-e...	7.258	3.1	ng/ul	106700	1	7.616	688298	20.0
1-Phenoxypropan...	11.163	3.2	ng/ul	168807	2	10.387	1048050	20.0
2,4,7,9-Tetrame...	13.169	6.7	ng/ul	508140	3	14.293	1515380	20.0
n-Hexadecanoic ...	18.092	27.0	ng/ul	2576850	4	17.134	1911640	20.0
Fluoran thene	19.281	3.5	ng/ul	339009	4	17.134	1911640	20.0
Octadecanoic acid	19.439	13.9	ng/ul	1258030	5	21.633	1803350	20.0
(DEL) Alkane: S...	21.310	6.0	ng/ul	537289	5	21.633	1803350	20.0