

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072721\
 Data File : BP006379.D
 Acq On : 28 Jul 2021 00:07
 Operator : CG/JU
 Sample : M3084-22
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0041

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: BP006379.D

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.305	32	37	45	rVB	95045	143706	1.38%	0.138%
2	3.575	80	83	97	rVB	25795	49249	0.47%	0.047%
3	3.705	101	105	129	rVB	779837	1326352	12.76%	1.273%
4	4.022	155	159	165	rVB	21721	34197	0.33%	0.033%
5	5.893	472	477	483	rBV	15110	25809	0.25%	0.025%
6	6.781	623	628	632	rBV2	12111	17223	0.17%	0.017%
7	6.957	651	658	669	rVB	1612939	2491778	23.97%	2.391%
8	7.128	680	687	698	rVB	1270704	1939879	18.66%	1.861%
9	7.316	711	719	727	rBV	1512512	2378418	22.88%	2.282%
10	7.416	732	736	744	rBV3	12424	26127	0.25%	0.025%
11	7.787	791	799	806	rBV	1438312	2302309	22.15%	2.209%
12	7.857	806	811	817	rVB4	10734	17870	0.17%	0.017%
13	8.493	910	919	928	rBV	2024766	3269293	31.45%	3.137%
14	8.940	987	995	1004	rBV	1491085	2394953	23.04%	2.298%
15	9.051	1010	1014	1019	rVB3	10093	15120	0.15%	0.015%
16	9.110	1020	1024	1028	rVB6	10233	13077	0.13%	0.013%
17	9.657	1109	1117	1134	rBV	1251343	2064676	19.86%	1.981%
18	9.887	1150	1156	1162	rVB	6366	12417	0.12%	0.012%
19	10.081	1182	1189	1197	rVB5	12262	24245	0.23%	0.023%
20	10.187	1197	1207	1217	rBV	1854634	3077054	29.60%	2.952%
21	10.357	1230	1236	1250	rBV	529570	801077	7.71%	0.769%
22	10.569	1264	1272	1280	rBV	2028257	3326371	32.00%	3.191%
23	10.710	1285	1296	1307	rBV	2045569	3481350	33.49%	3.340%
24	12.075	1523	1528	1532	rBV6	7572	14474	0.14%	0.014%
25	12.151	1536	1541	1551	rVB	34886	59800	0.58%	0.057%
26	13.222	1716	1723	1726	rVV5	11684	25908	0.25%	0.025%
27	13.257	1726	1729	1733	rVV5	10042	15953	0.15%	0.015%
28	13.322	1735	1740	1746	rVV	63662	104572	1.01%	0.100%
29	13.404	1750	1754	1761	rVB2	14831	28252	0.27%	0.027%
30	13.834	1820	1827	1836	rBV	3378763	4716356	45.37%	4.525%
31	14.104	1862	1873	1885	rBV	3092167	4721214	45.42%	4.530%
32	14.251	1892	1898	1908	rVV2	36639	88398	0.85%	0.085%
33	14.322	1908	1910	1917	rVB8	10488	14019	0.13%	0.013%
34	14.410	1917	1925	1938	rVB	2937434	4142003	39.85%	3.974%
35	14.610	1949	1959	1977	rBV	1940731	2908000	27.98%	2.790%
36	14.839	1994	1998	2004	rVB2	19931	28179	0.27%	0.027%

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 Title : SVOA CALIBRATION

37	14.957	2011	2018	2024	rVB	22993	43887	0.42%	0.042%
38	15.222	2060	2063	2068	rBV4	5594	10871	0.10%	0.010%
39	15.281	2069	2073	2078	rVB4	11144	13799	0.13%	0.013%
40	15.404	2087	2094	2105	rVB	4255715	6170352	59.36%	5.920%
41	15.522	2107	2114	2127	rBV	1646245	2210915	21.27%	2.121%
42	15.645	2130	2135	2141	rVB2	50884	73538	0.71%	0.071%
43	15.975	2186	2191	2194	rBV6	13286	21079	0.20%	0.020%
44	16.092	2206	2211	2216	rBV8	15072	29401	0.28%	0.028%
45	16.204	2225	2230	2236	rVB6	21333	34440	0.33%	0.033%
46	16.310	2242	2248	2255	rBV9	16713	30889	0.30%	0.030%
47	16.404	2260	2264	2267	rBV2	14575	23772	0.23%	0.023%
48	16.563	2287	2291	2296	rVB3	27211	41556	0.40%	0.040%
49	16.675	2307	2310	2314	rVB3	10152	12444	0.12%	0.012%
50	16.910	2345	2350	2352	rBV5	16020	23109	0.22%	0.022%
51	16.939	2352	2355	2359	rVB3	23370	31494	0.30%	0.030%
52	17.163	2386	2393	2402	rBV	3183422	4541066	43.69%	4.357%
53	17.263	2403	2410	2417	rBV	5834564	7873980	75.75%	7.554%
54	17.345	2422	2424	2434	rVB	10190	23182	0.22%	0.022%
55	17.539	2453	2457	2458	rBV2	10633	12469	0.12%	0.012%
56	17.716	2484	2487	2493	rVB7	10347	16906	0.16%	0.016%
57	17.851	2506	2510	2516	rVB	160022	196961	1.89%	0.189%
58	18.069	2542	2547	2556	rBV	921890	1280375	12.32%	1.228%
59	18.363	2593	2597	2605	rBV	14350	29903	0.29%	0.029%
60	18.963	2695	2699	2700	rBV2	59942	73352	0.71%	0.070%
61	18.986	2700	2703	2708	rVB	441077	522837	5.03%	0.502%
62	19.080	2715	2719	2722	rBV2	87275	112874	1.09%	0.108%
63	19.121	2722	2726	2728	rVV	109982	126531	1.22%	0.121%
64	19.157	2728	2732	2735	rVV2	100191	187621	1.81%	0.180%
65	19.186	2735	2737	2745	rVB5	49050	79131	0.76%	0.076%
66	19.304	2753	2757	2768	rBV	565517	795068	7.65%	0.763%
67	19.457	2780	2783	2785	rBV	27471	27278	0.26%	0.026%
68	19.504	2785	2791	2799	rVV	8150330	10394512	100.00%	9.973%
69	19.569	2799	2802	2806	rVB2	94830	107313	1.03%	0.103%
70	20.016	2875	2878	2881	rBV4	35492	47973	0.46%	0.046%
71	20.051	2881	2884	2887	rVV	157127	156869	1.51%	0.151%
72	20.539	2962	2967	2970	rBV2	255300	321784	3.10%	0.309%
73	20.574	2970	2973	2974	rBV3	48102	50274	0.48%	0.048%
74	20.986	3039	3043	3050	rBV2	1264642	1614875	15.54%	1.549%
75	21.051	3052	3054	3059	rVB	90444	106602	1.03%	0.102%
76	21.257	3084	3089	3097	rBV2	3457800	4421770	42.54%	4.242%

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

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Peak separation: 5

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Title : SVOA CALIBRATION

77	21.427	3114	3118	3122	rBV	406485	459639	4.42%	0.441%
78	21.880	3192	3195	3200	rBV	403309	517186	4.98%	0.496%
79	22.368	3275	3278	3290	rVB	471297	843331	8.11%	0.809%
80	22.915	3367	3371	3377	rVB2	407834	623278	6.00%	0.598%
81	23.457	3455	3463	3471	rVV2	4455910	8814850	84.80%	8.457%
82	23.533	3472	3476	3481	rVV	322073	569787	5.48%	0.547%
83	23.604	3482	3488	3501	rVB2	2001963	3851556	37.05%	3.695%
84	24.245	3593	3597	3604	rVB	296079	558660	5.37%	0.536%

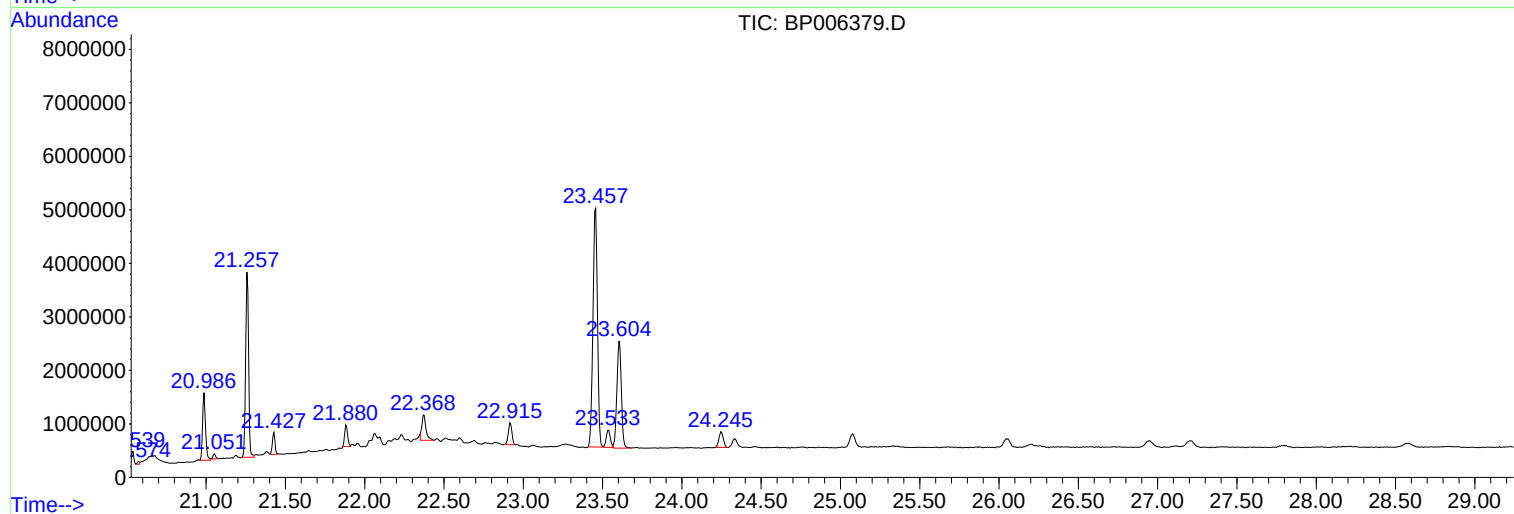
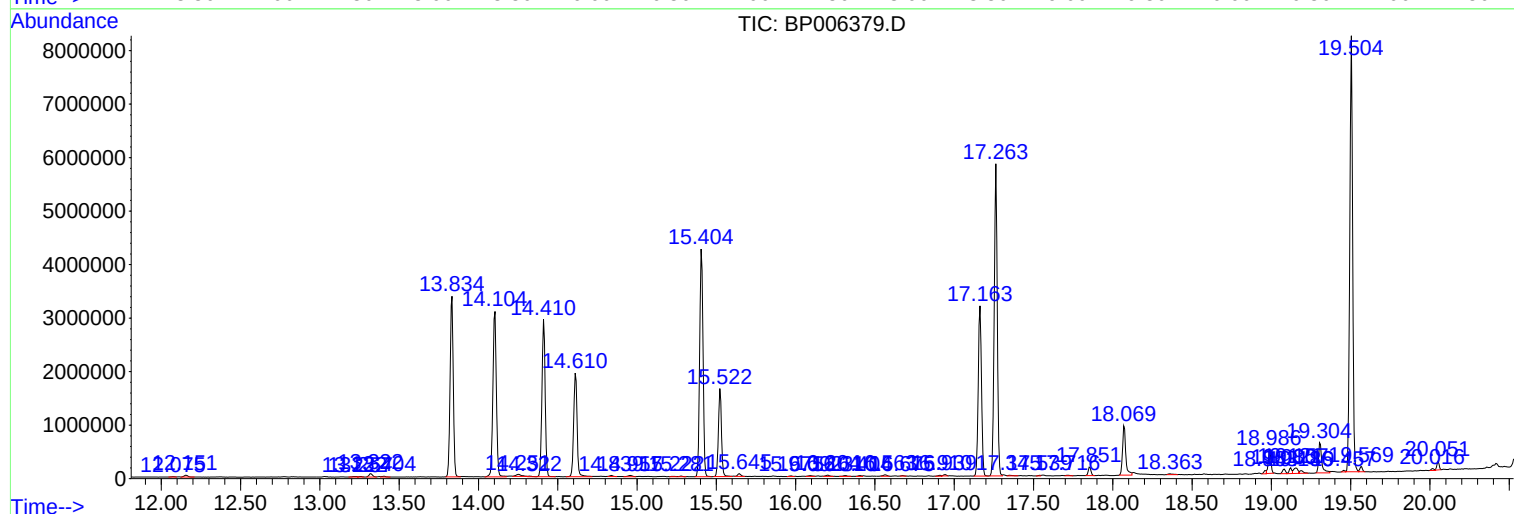
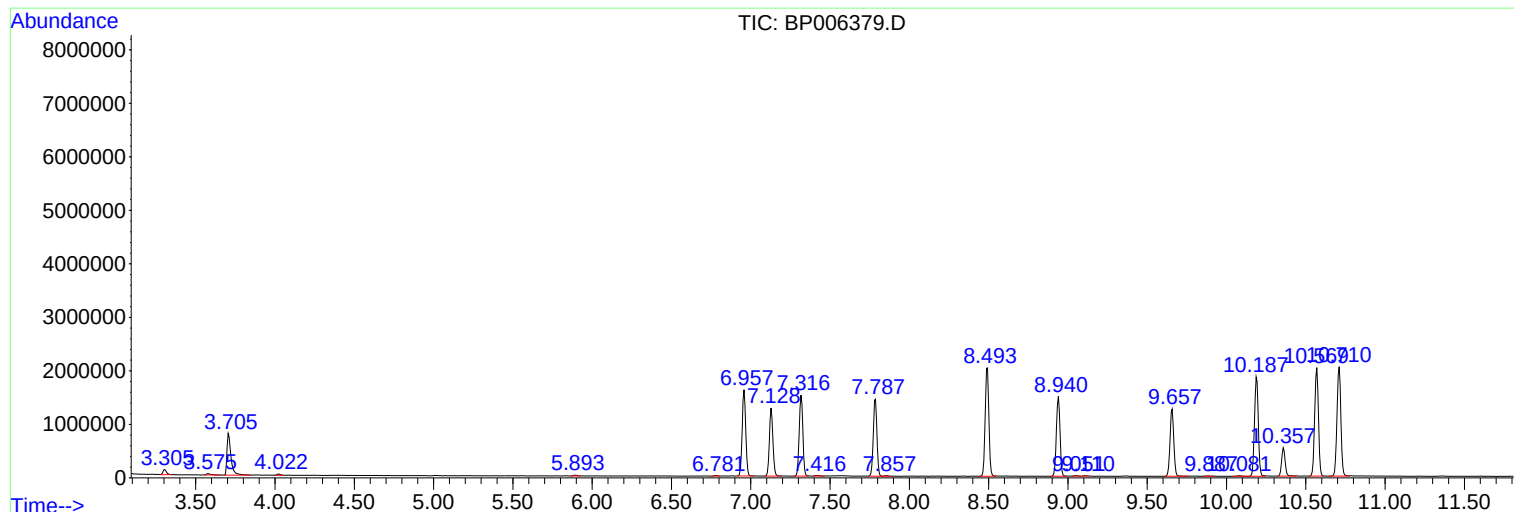
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP072421.M
Quant Title : SVOA CALIBRATION

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



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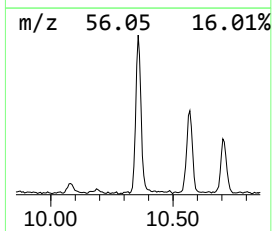
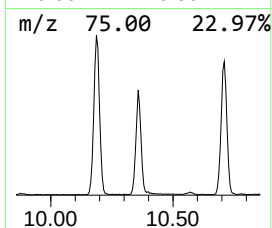
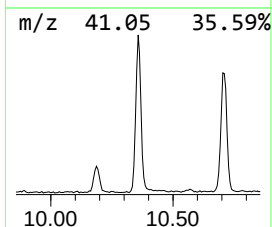
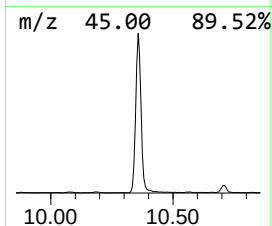
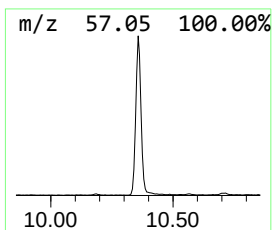
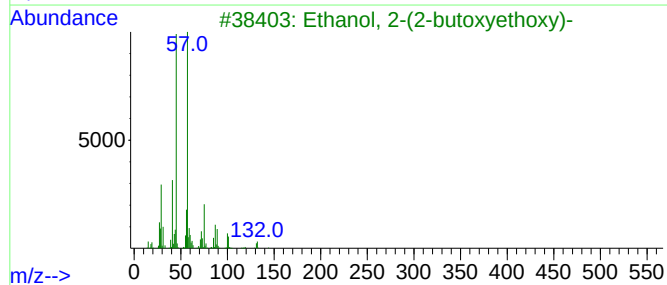
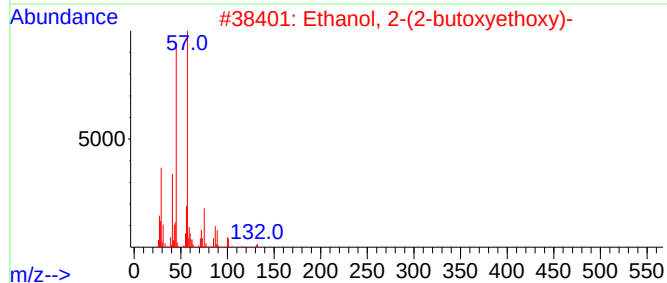
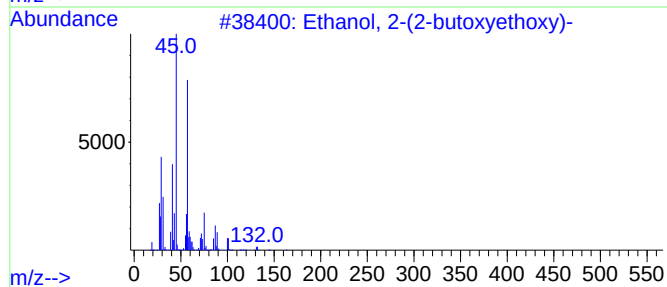
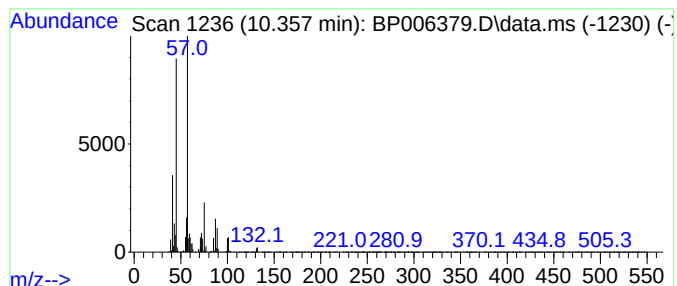
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP072421.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Ethanol, 2-(2-butoxyethoxy)- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.360	4.82 ng/ul	801077	Naphthalene-d8	10.569

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	86
4			Ethanol, 1-(2-butoxyethoxy)-	162	C8H18O3	054446-78-5	83
5			Ethanol, 1-(2-butoxyethoxy)-	162	C8H18O3	054446-78-5	83



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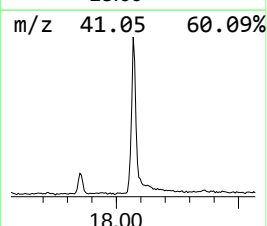
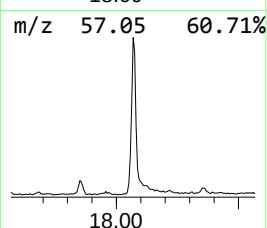
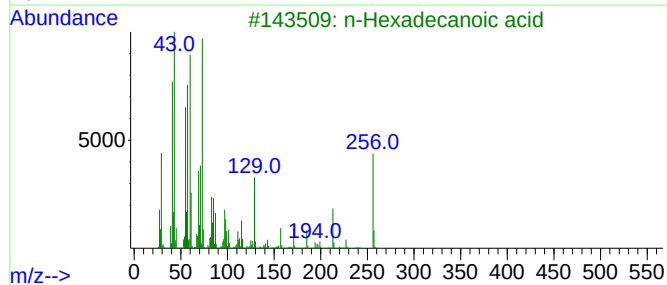
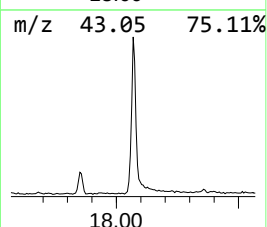
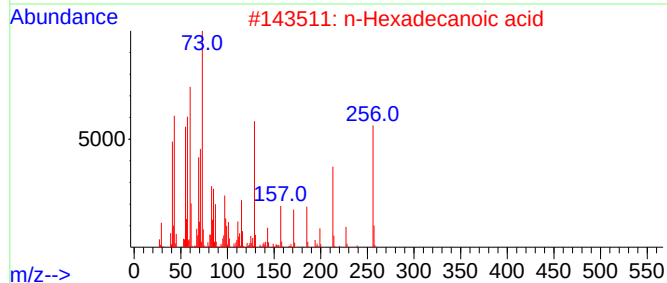
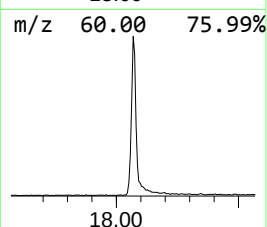
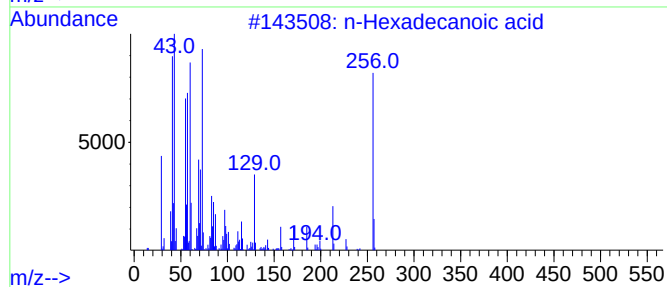
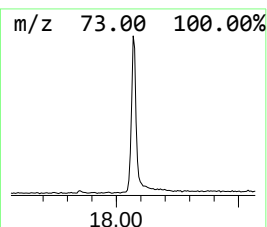
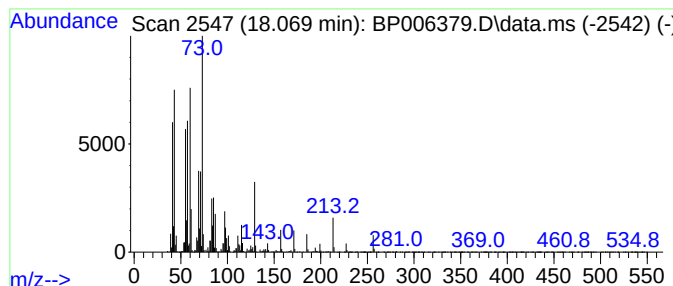
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP072421.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 n-Hexadecanoic acid Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.070	5.64 ng/ul	1280380	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	99
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	95



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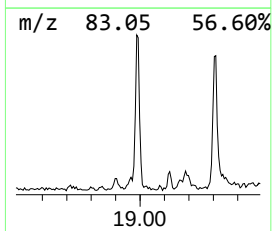
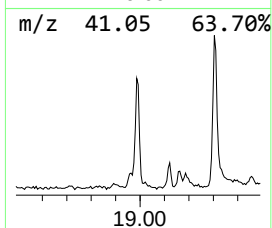
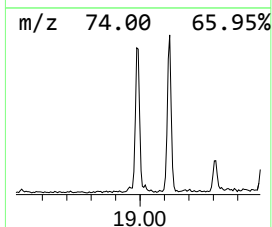
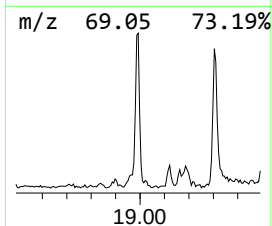
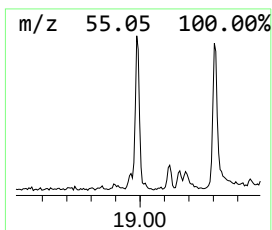
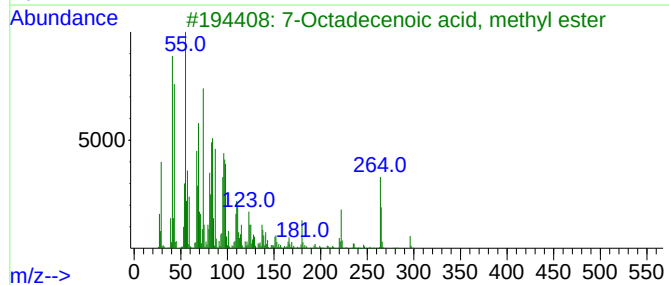
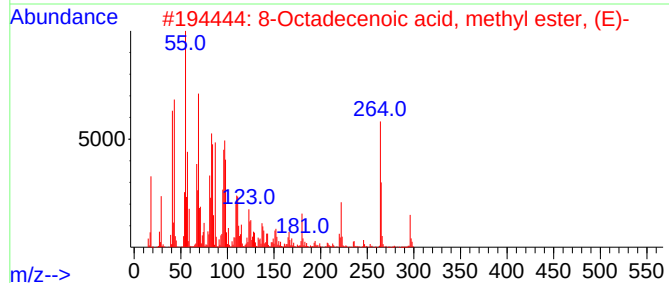
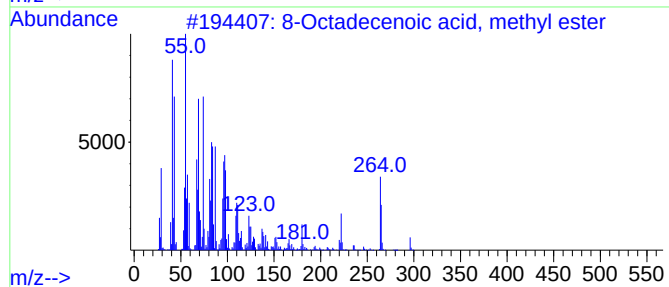
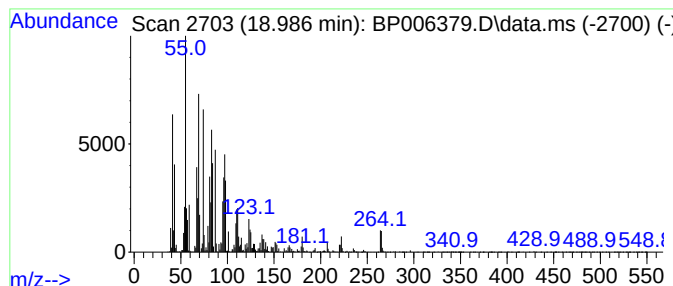
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP072421.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 8-Octadecenoic acid, methyl... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.990	2.30 ng/ul	522837	Phenanthrene-d10	17.163

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			8-Octadecenoic acid, methyl ester	296	C19H36O2	002345-29-1	99
2			8-Octadecenoic acid, methyl este...	296	C19H36O2	026528-50-7	99
3			7-Octadecenoic acid, methyl ester	296	C19H36O2	057396-98-2	97
4			9-Octadecenoic acid (Z)-, methyl...	296	C19H36O2	000112-62-9	97
5			(Z)-15-Octadecenoic acid methyl ...	296	C19H36O2	1000427-69-3	96



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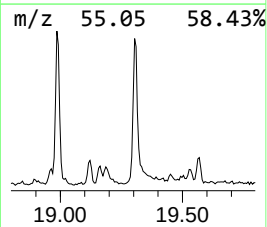
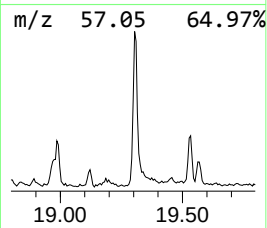
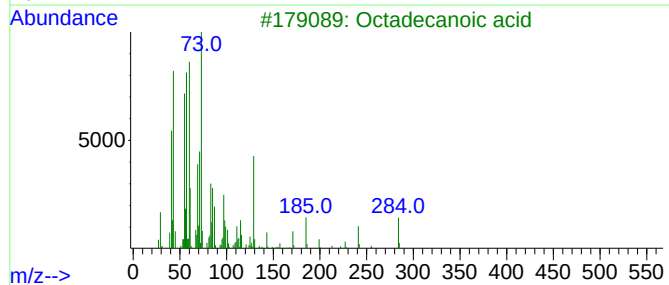
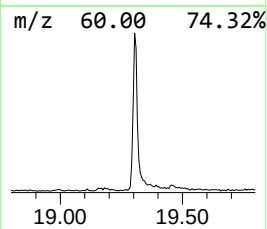
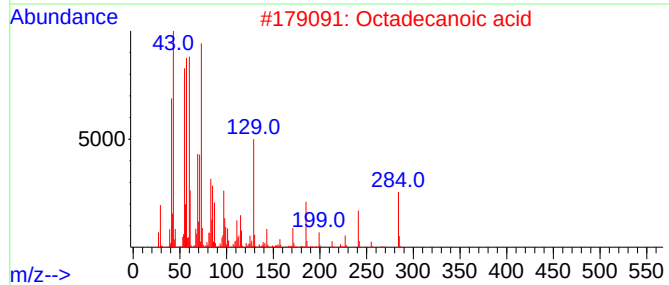
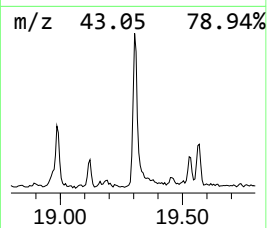
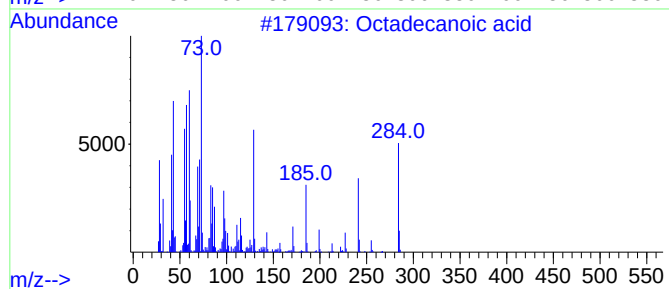
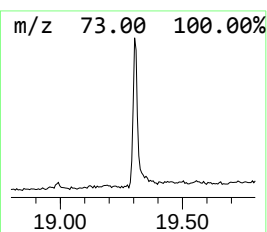
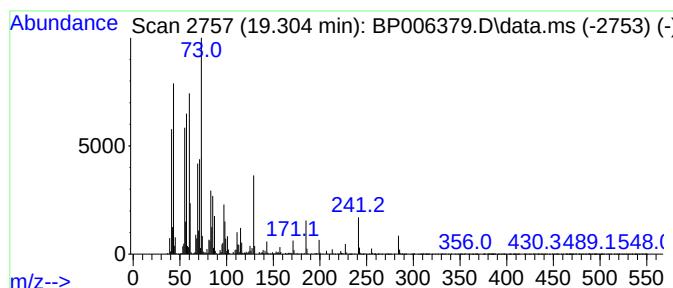
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 Octadecanoic acid Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.300	3.60 ng/ul	795068	Chrysene-d12	21.257

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	98
4			Octadecanoic acid	284	C18H36O2	000057-11-4	96
5			Octadecanoic acid	284	C18H36O2	000057-11-4	95



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072721\
Data File : BP006379.D
Acq On : 28 Jul 2021 00:07
Operator : CG/JU
Sample : M3084-22
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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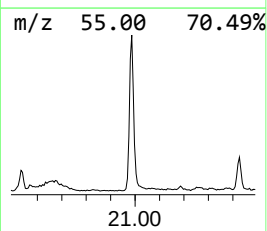
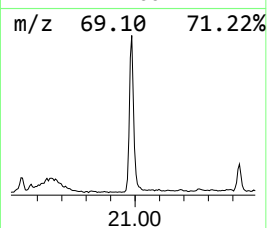
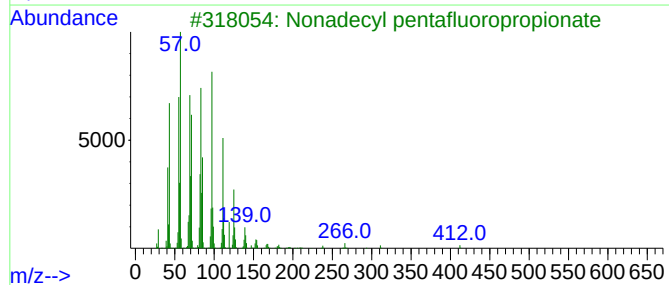
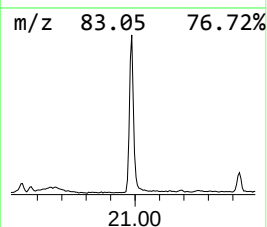
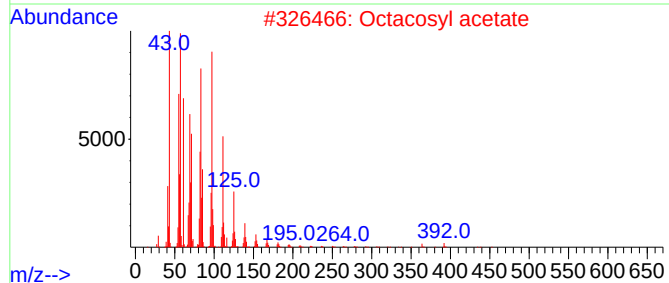
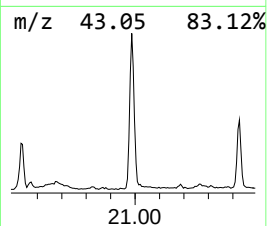
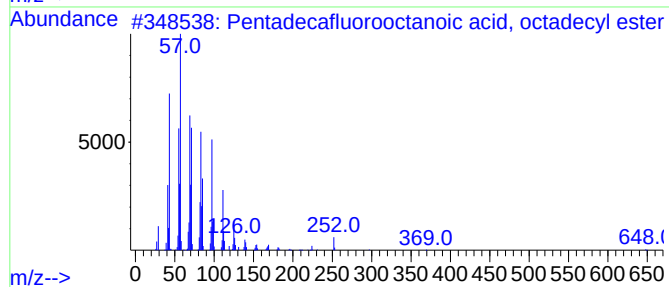
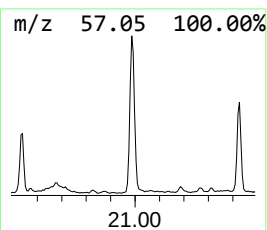
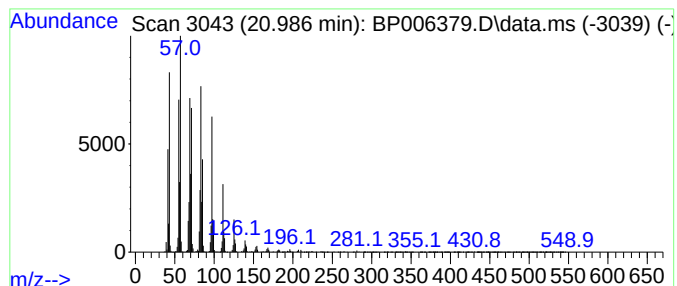
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 Pentadecafluorooctanoic aci... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.990	7.30 ng/ul	1614880	Chrysene-d12	21.257

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentadecafluorooctanoic acid, oc...	666	C26H37F15O2	1000406-04-8	94
2			Octacosyl acetate	452	C30H60O2	018206-97-8	93
3			Nonadecyl pentafluoropropionate	430	C22H39F5O2	1000351-88-8	91
4			Ethanol, 2-(tetradecyloxy)-	258	C16H34O2	002136-70-1	91
5			Pentafluoropropionic acid, octad...	416	C21H37F5O2	959261-25-7	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072721\
Data File : BP006379.D
Acq On : 28 Jul 2021 00:07
Operator : CG/JU
Sample : M3084-22
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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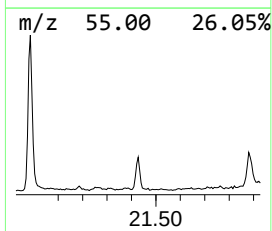
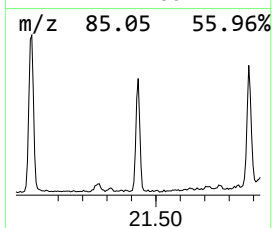
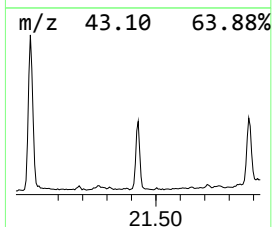
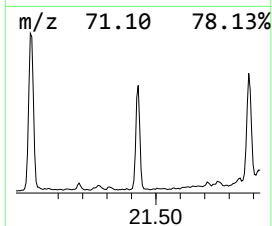
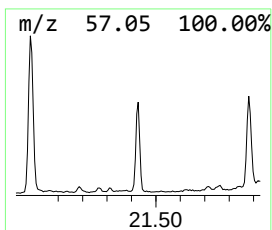
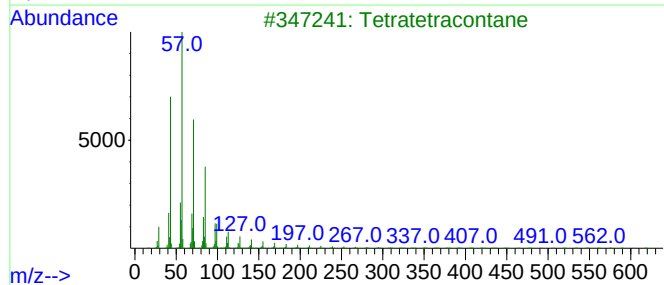
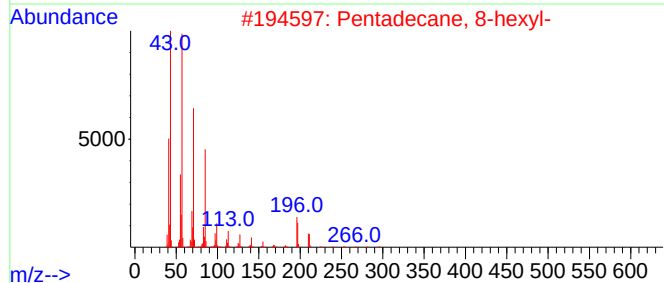
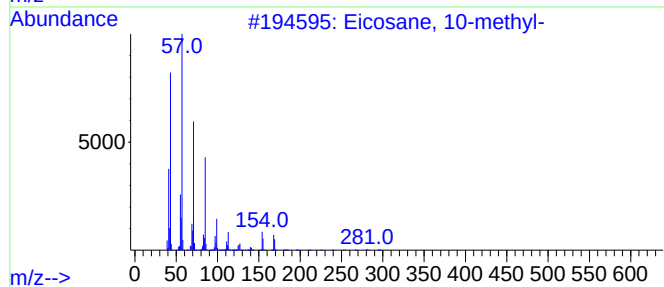
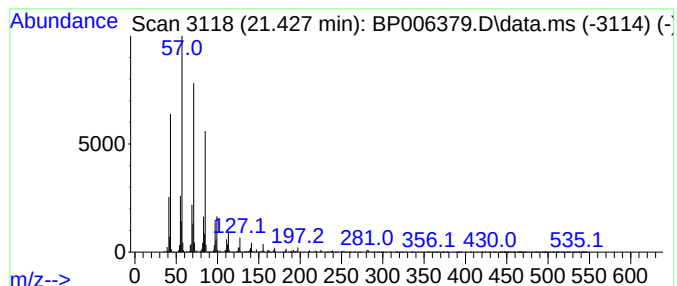
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 (DEL) Alkane: Straight-Chai... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.430	2.08 ng/ul	459639	Chrysene-d12	21.257

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Eicosane, 10-methyl-	296	C21H44	054833-23-7	93
2			Pentadecane, 8-hexyl-	296	C21H44	013475-75-7	93
3			Tetratetracontane	619	C44H90	007098-22-8	91
4			Pentacosane	352	C25H52	000629-99-2	91
5			Hexatriacontane	507	C36H74	000630-06-8	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072721\
Data File : BP006379.D
Acq On : 28 Jul 2021 00:07
Operator : CG/JU
Sample : M3084-22
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampled :
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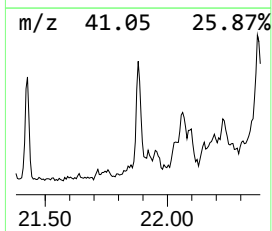
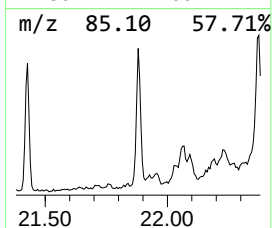
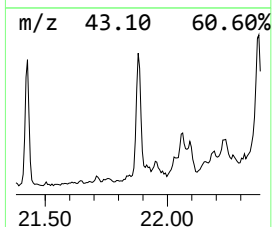
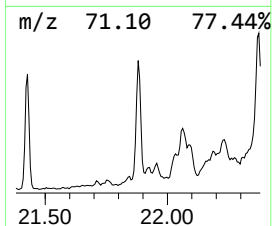
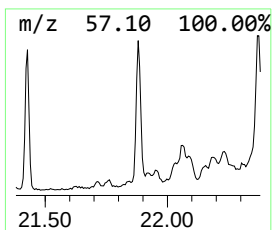
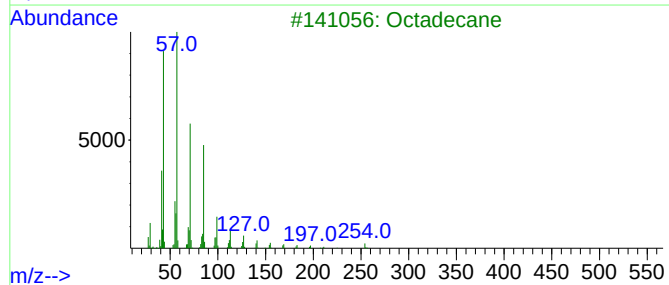
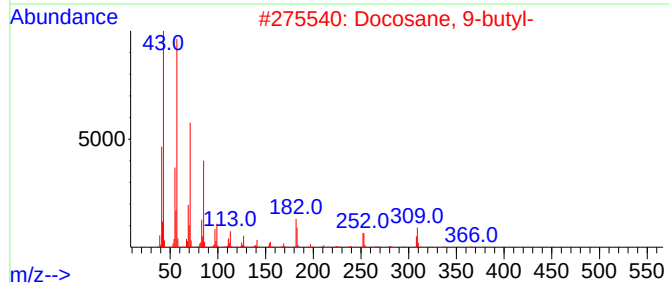
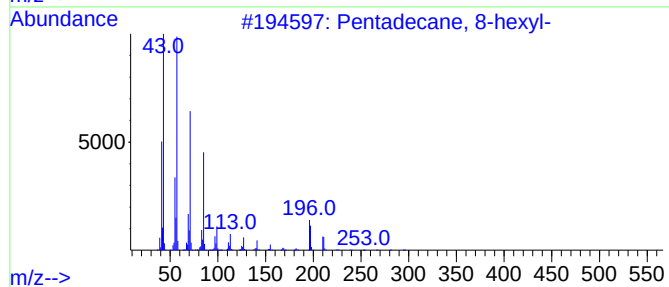
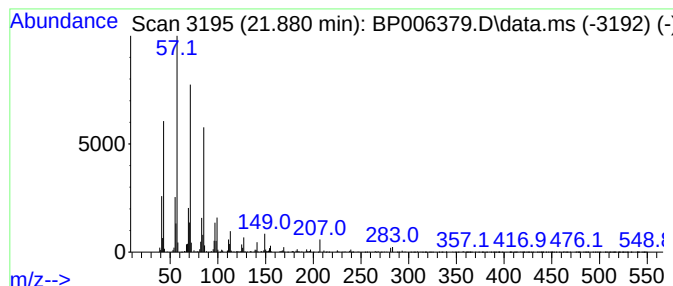
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 (DEL) Alkane: Straight-Chai... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.880	2.34 ng/ul	517186	Chrysene-d12	21.257

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentadecane, 8-hexyl-	296	C21H44	013475-75-7	93
2			Docosane, 9-butyl-	366	C26H54	055282-14-9	93
3			Octadecane	254	C18H38	000593-45-3	92
4			Octadecane, 5,14-dibutyl-	366	C26H54	055282-13-8	90
5			Tetracosane, 3-ethyl-	366	C26H54	055282-17-2	89



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072721\
Data File : BP006379.D
Acq On : 28 Jul 2021 00:07
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Sample : M3084-22
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument :
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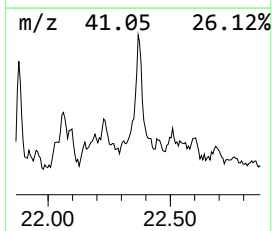
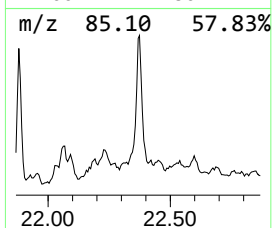
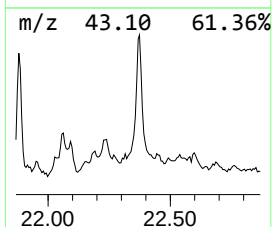
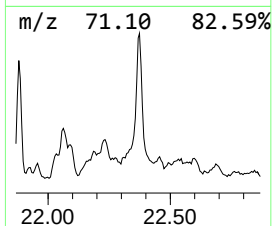
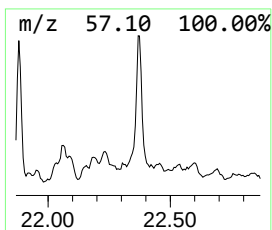
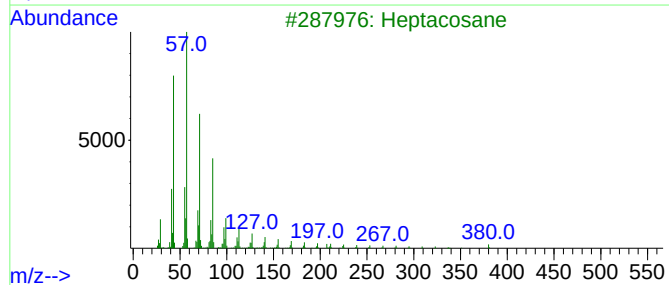
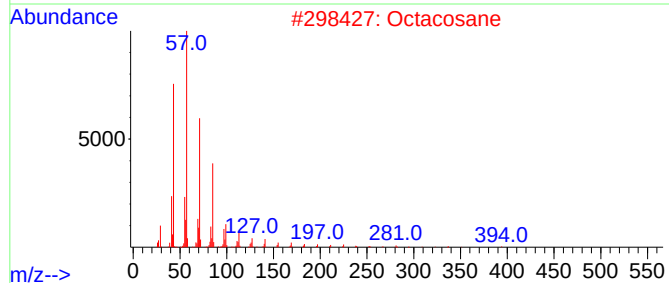
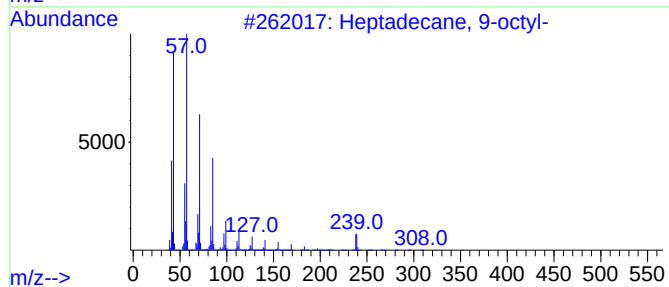
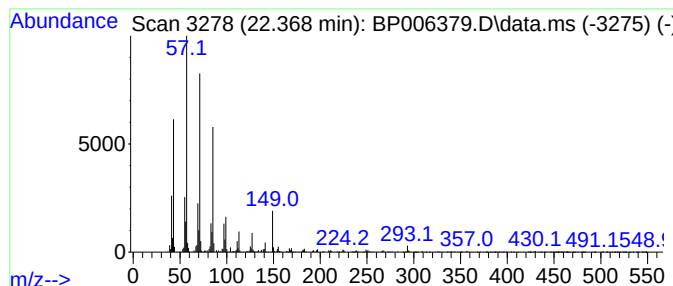
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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.370	3.81 ng/ul	843331	Chrysene-d12	21.257

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heptadecane, 9-octyl-	352	C25H52	007225-64-1	87
2			Octacosane	394	C28H58	000630-02-4	87
3			Heptacosane	380	C27H56	000593-49-7	83
4			Nonadecane	268	C19H40	000629-92-5	83
5			Hexacosane	366	C26H54	000630-01-3	83



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072721\
Data File : BP006379.D
Acq On : 28 Jul 2021 00:07
Operator : CG/JU
Sample : M3084-22
Misc :
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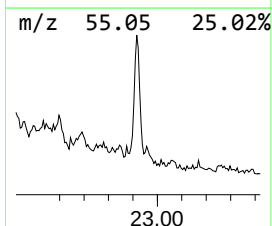
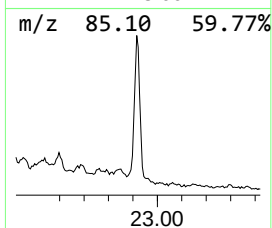
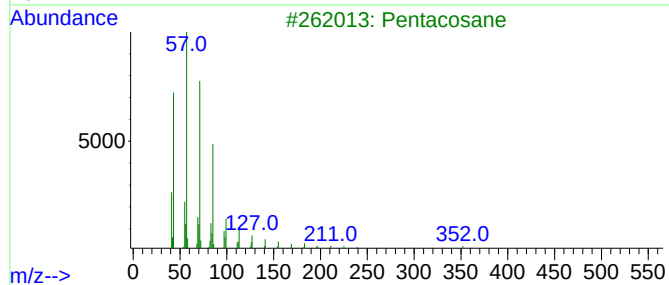
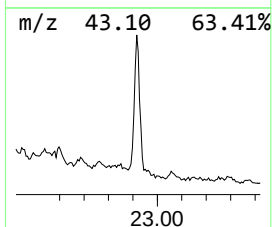
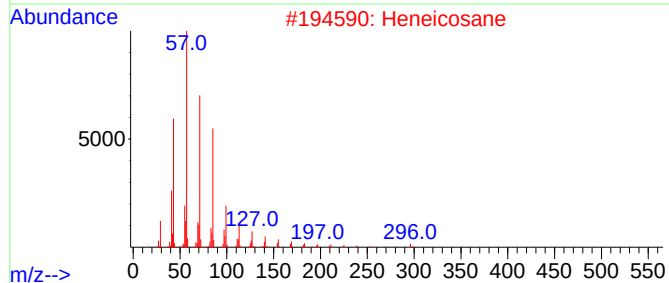
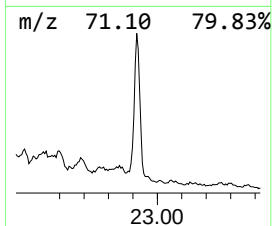
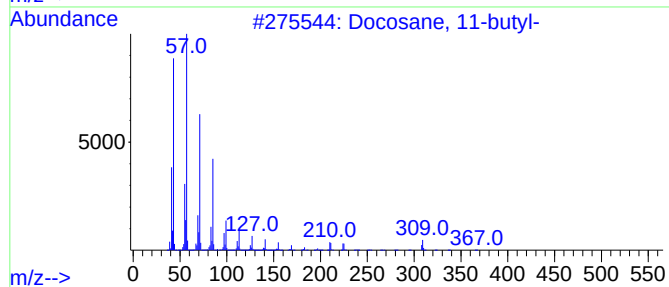
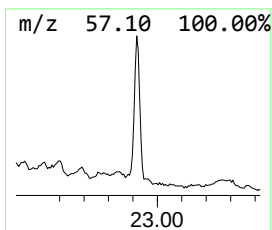
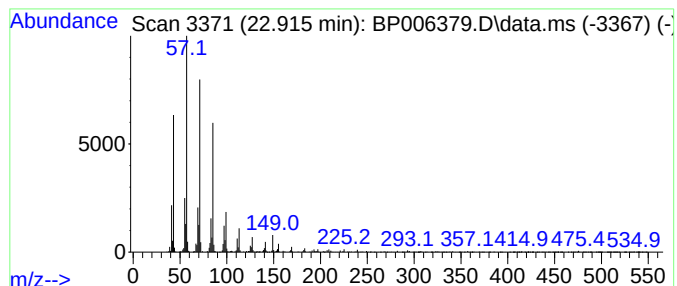
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 (DEL) Alkane: Straight-Chai... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD		R.T.	
22.920	3.24 ng/ul	623278	Perylene-d12		23.604	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Docosane, 11-butyl-		366	C26H54	013475-76-8	94
2	Heneicosane		296	C21H44	000629-94-7	91
3	Pentacosane		352	C25H52	000629-99-2	91
4	Octadecane, 5,14-dibutyl-		366	C26H54	055282-13-8	90
5	Heneicosane, 11-pentyl-		366	C26H54	014739-72-1	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072721\
Data File : BP006379.D
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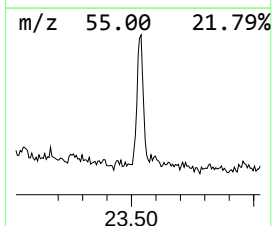
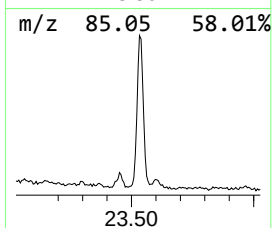
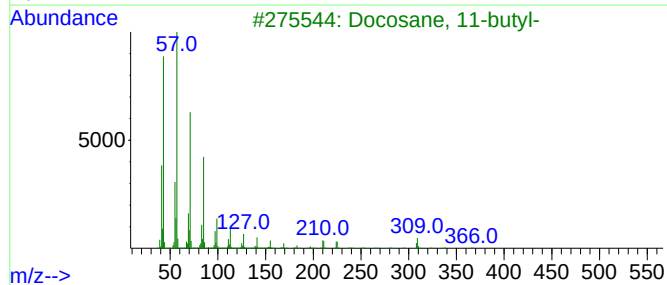
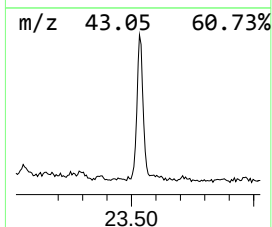
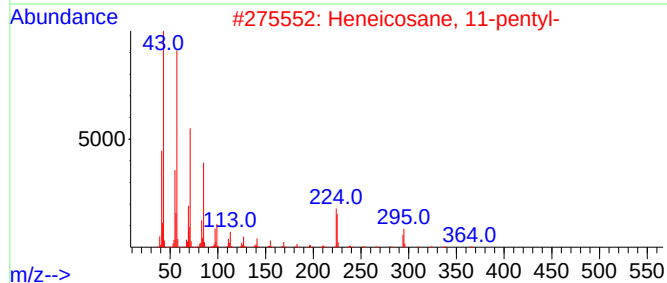
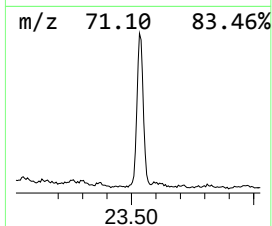
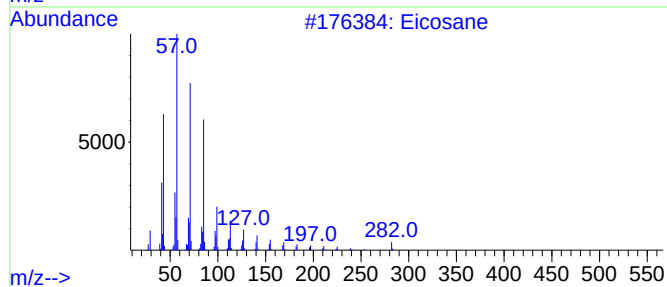
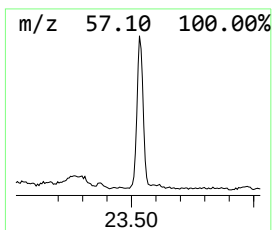
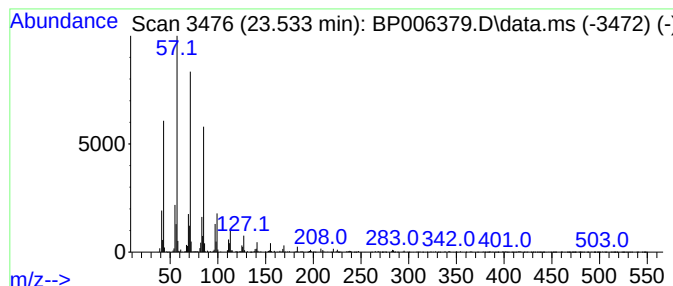
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 (DEL) Alkane: Straight-Chai... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.530	2.96 ng/ul	569787	Perylene-d12	23.604

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Eicosane			282	C20H42	000112-95-8	95
2	Heneicosane, 11-pentyl-			366	C26H54	014739-72-1	94
3	Docosane, 11-butyl-			366	C26H54	013475-76-8	94
4	Nonadecane			268	C19H40	000629-92-5	93
5	Heneicosane			296	C21H44	000629-94-7	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072721\
Data File : BP006379.D
Acq On : 28 Jul 2021 00:07
Operator : CG/JU
Sample : M3084-22
Misc :
ALS Vial : 24 Sample Multiplier: 1

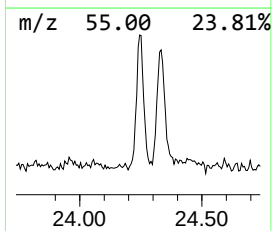
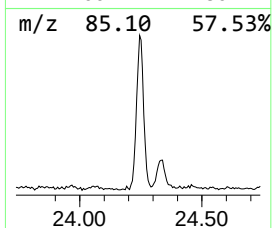
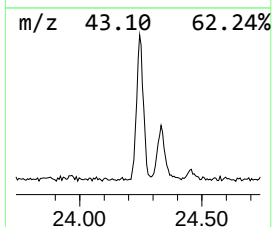
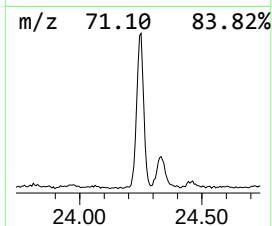
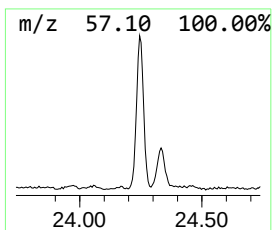
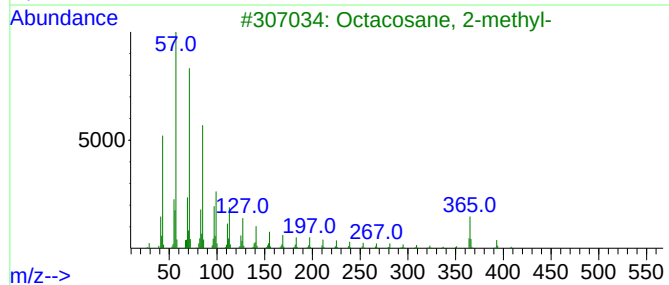
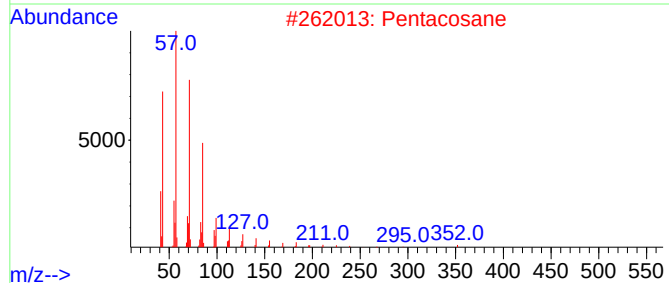
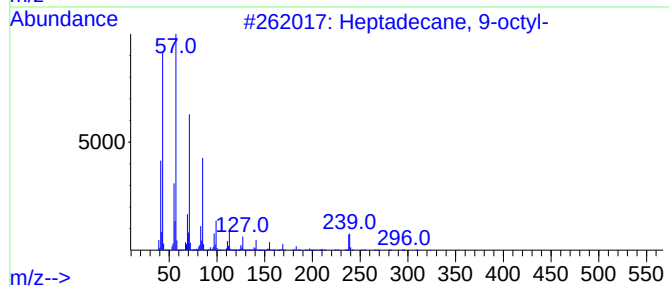
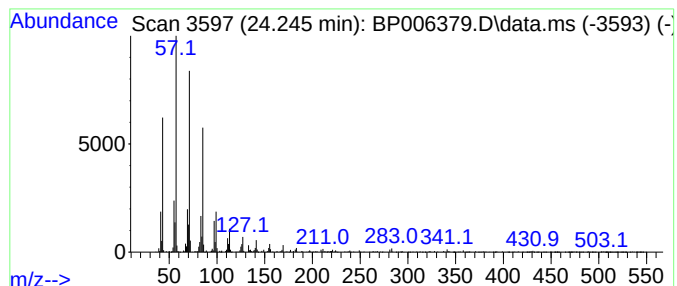
Instrument :
BNA_P
ClientSampleId :
C0041

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 11 (DEL) Alkane: Straight-Chai... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.	
24.240	2.90 ng/ul	558660	Perylene-d12	23.604	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptadecane, 9-octyl-	352	C25H52	007225-64-1	95
2	Pentacosane	352	C25H52	000629-99-2	91
3	Octacosane, 2-methyl-	408	C29H60	001560-98-1	91
4	Octacosane, 2-methyl-	408	C29H60	001560-98-1	91
5	Heneicosane	296	C21H44	000629-94-7	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072721\
Data File : BP006379.D
Acq On : 28 Jul 2021 00:07
Operator : CG/JU
Sample : M3084-22
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0041

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-(2-b...	10.360	4.8	ng/ul	801077	2	10.569	3326370	20.0
n-Hexadecanoic ...	18.070	5.6	ng/ul	1280380	4	17.163	4541070	20.0
8-Octadecenoic ...	18.990	2.3	ng/ul	522837	4	17.163	4541070	20.0
Octadecanoic acid	19.300	3.6	ng/ul	795068	5	21.257	4421770	20.0
Pentadecafluoro...	20.990	7.3	ng/ul	1614880	5	21.257	4421770	20.0
(DEL) Alkane: S...	21.430	2.1	ng/ul	459639	5	21.257	4421770	20.0
(DEL) Alkane: S...	21.880	2.3	ng/ul	517186	5	21.257	4421770	20.0
(DEL) Alkane: S...	22.370	3.8	ng/ul	843331	5	21.257	4421770	20.0
(DEL) Alkane: S...	22.920	3.2	ng/ul	623278	6	23.604	3851560	20.0
(DEL) Alkane: S...	23.530	3.0	ng/ul	569787	6	23.604	3851560	20.0
(DEL) Alkane: S...	24.240	2.9	ng/ul	558660	6	23.604	3851560	20.0