

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110722\
 Data File : BP012536.D
 Acq On : 07 Nov 2022 17:57
 Operator : CG/JU
 Sample : N5353-16
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 DBWX1

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-BP102522.M
 Title : SVOA CALIBRATION

Signal : TIC: BP012536.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.234	38	42	45	rBV	43086	58688	0.63%	0.069%
2	3.663	112	115	136	rBV	173785	373641	3.98%	0.442%
3	3.863	144	149	155	rVB2	18842	31783	0.34%	0.038%
4	3.958	162	165	171	rVB	14006	22695	0.24%	0.027%
5	5.093	352	358	369	rVB	69829	122338	1.30%	0.145%
6	6.957	669	675	688	rBV	212352	428943	4.57%	0.507%
7	7.116	695	702	717	rBV	936523	1528849	16.30%	1.809%
8	7.310	728	735	751	rBV	842998	1426800	15.21%	1.688%
9	7.446	753	758	765	rVB10	7539	14740	0.16%	0.017%
10	7.775	806	814	825	rBV	870256	1467980	15.65%	1.737%
11	8.134	870	875	881	rVB5	14511	29289	0.31%	0.035%
12	8.499	929	937	956	rBV	700290	1280681	13.65%	1.515%
13	8.946	1006	1013	1027	rBV	1165526	1973509	21.04%	2.335%
14	9.157	1045	1049	1054	rVB8	7962	13299	0.14%	0.016%
15	9.293	1066	1072	1074	rBV6	12781	22665	0.24%	0.027%
16	9.322	1074	1077	1086	rVB2	26796	47647	0.51%	0.056%
17	9.604	1118	1125	1129	rBV4	9337	18807	0.20%	0.022%
18	9.669	1129	1136	1148	rVV	850651	1504919	16.04%	1.780%
19	9.951	1177	1184	1188	rBV10	7146	17797	0.19%	0.021%
20	10.063	1199	1203	1209	rVB6	6493	9866	0.11%	0.012%
21	10.204	1220	1227	1248	rBV	1328972	2574895	27.45%	3.046%
22	10.363	1248	1254	1280	rVV	461637	1178819	12.56%	1.395%
23	10.575	1283	1290	1306	rVV	1408088	2417796	25.77%	2.860%
24	10.728	1309	1316	1330	rVV	1088203	2071726	22.08%	2.451%
25	10.840	1330	1335	1344	rVB2	43426	89449	0.95%	0.106%
26	11.440	1434	1437	1441	rVB6	6498	10637	0.11%	0.013%
27	11.616	1462	1467	1471	rBV7	7166	14862	0.16%	0.018%
28	11.834	1493	1504	1511	rVB	45310	90800	0.97%	0.107%
29	12.034	1535	1538	1543	rVB2	10314	13900	0.15%	0.016%
30	12.175	1554	1562	1571	rVV3	18959	47768	0.51%	0.057%
31	12.251	1571	1575	1583	rVB6	16712	34587	0.37%	0.041%
32	12.328	1583	1588	1595	rBV8	10238	19428	0.21%	0.023%
33	12.387	1596	1598	1608	rVB10	4686	11227	0.12%	0.013%
34	12.498	1608	1617	1623	rBV4	18024	45151	0.48%	0.053%
35	12.804	1667	1669	1673	rVB5	10636	12064	0.13%	0.014%
36	12.951	1687	1694	1700	rBV3	28985	57781	0.62%	0.068%

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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-BP102522.M
 Title : SVOA CALIBRATION

37	13.034	1704	1708	1713	rBV6	10414	20563	0.22%	0.024%
38	13.151	1725	1728	1731	rBV5	8881	12124	0.13%	0.014%
39	13.239	1739	1743	1746	rVB2	16133	20548	0.22%	0.024%
40	13.322	1751	1757	1765	rBV	461956	717409	7.65%	0.849%
41	13.392	1765	1769	1772	rVV6	13476	24120	0.26%	0.029%
42	13.428	1772	1775	1780	rVB6	17109	28722	0.31%	0.034%
43	13.745	1826	1829	1832	rBV5	7443	10958	0.12%	0.013%
44	13.845	1837	1846	1868	rBV	2973561	4469838	47.64%	5.288%
45	14.122	1886	1893	1909	rBV	3374640	5105726	54.42%	6.040%
46	14.316	1924	1926	1931	rVB6	10893	11292	0.12%	0.013%
47	14.428	1938	1945	1960	rBV	2341568	3548222	37.82%	4.197%
48	14.681	1983	1988	2002	rBV2	91186	273045	2.91%	0.323%
49	14.792	2002	2007	2011	rVV	38195	55263	0.59%	0.065%
50	14.851	2011	2017	2022	rVV5	25133	59595	0.64%	0.070%
51	14.969	2034	2037	2043	rVB3	35031	53970	0.58%	0.064%
52	15.039	2045	2049	2054	rBV8	13295	25454	0.27%	0.030%
53	15.181	2070	2073	2078	rBV7	10002	17885	0.19%	0.021%
54	15.275	2085	2089	2094	rVB3	21777	32422	0.35%	0.038%
55	15.428	2108	2115	2131	rBV	4569279	6906914	73.62%	8.171%
56	15.563	2132	2138	2152	rBV	1271725	1871664	19.95%	2.214%
57	15.881	2189	2192	2199	rVB7	32680	50570	0.54%	0.060%
58	17.192	2408	2415	2425	rBV	2972154	4481904	47.77%	5.302%
59	17.292	2425	2432	2444	rVV	5354570	7810725	83.25%	9.240%
60	17.792	2515	2517	2521	rVB2	20555	23240	0.25%	0.027%
61	17.857	2524	2528	2537	rBV	255270	325898	3.47%	0.386%
62	18.080	2561	2566	2575	rBV	409133	673190	7.18%	0.796%
63	18.963	2714	2716	2717	rBV2	39511	33392	0.36%	0.040%
64	18.992	2717	2721	2735	rVB	1017749	1257054	13.40%	1.487%
65	19.122	2737	2743	2748	rBV2	399671	529832	5.65%	0.627%
66	19.180	2749	2753	2756	rBV	82081	116306	1.24%	0.138%
67	19.316	2772	2776	2787	rBV3	290927	506224	5.40%	0.599%
68	19.533	2807	2813	2825	rBV	7138585	9381868	100.00%	11.099%
69	20.374	2954	2956	2961	rVB5	77171	86068	0.92%	0.102%
70	21.004	3060	3063	3071	rVB4	206732	333025	3.55%	0.394%
71	21.286	3106	3111	3119	rBV	3486217	4586363	48.89%	5.426%
72	23.515	3483	3490	3503	rVB2	3887503	7935441	84.58%	9.387%
73	23.668	3510	3516	3528	rVB2	2047544	4049532	43.16%	4.791%

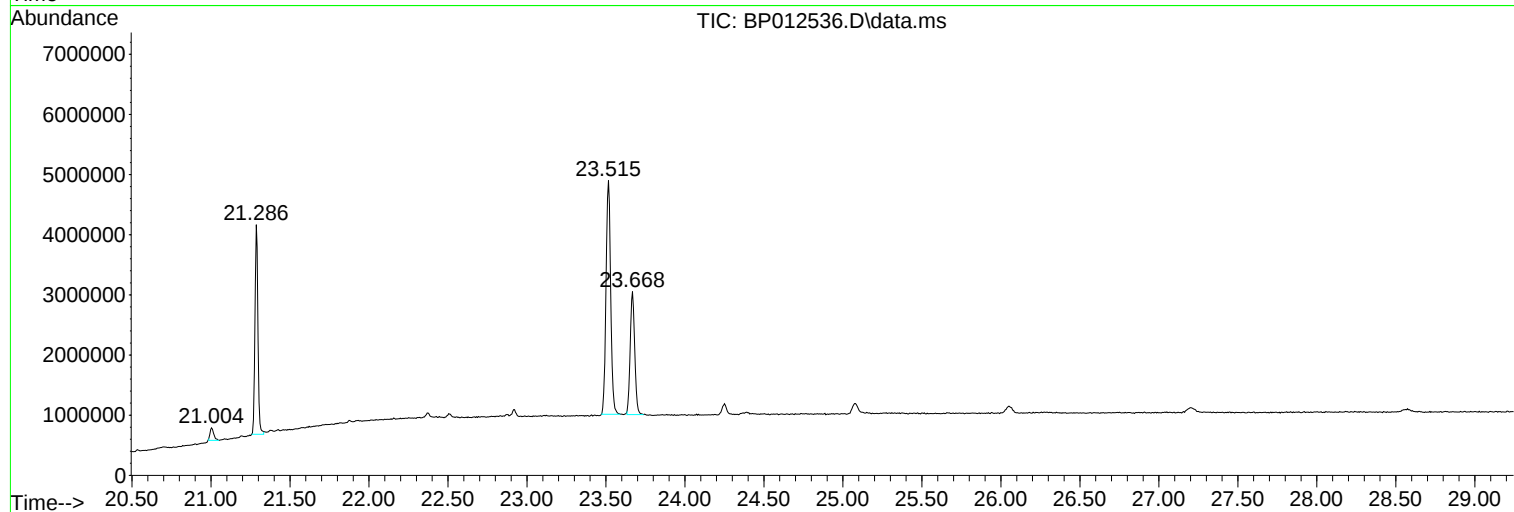
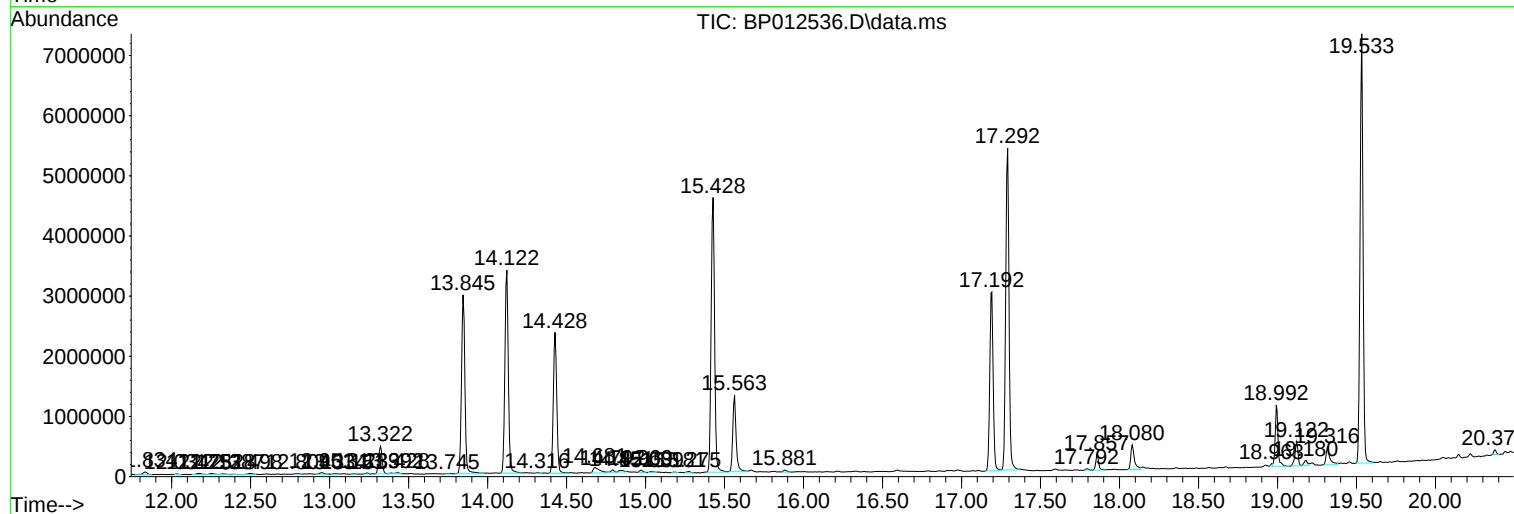
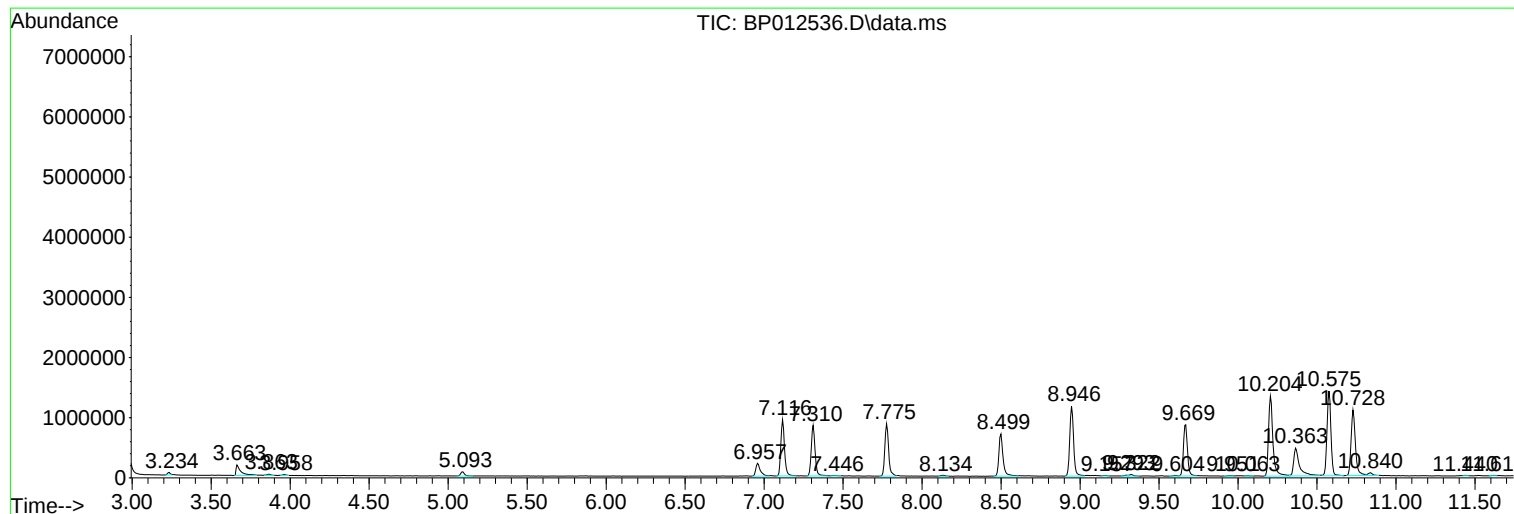
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
Quant Title : SVOA CALIBRATION

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



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Sample : N5353-16
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ALS Vial : 13 Sample Multiplier: 1

Instrument :
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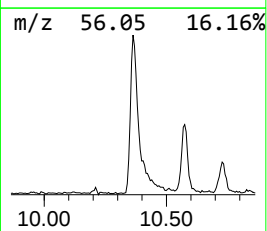
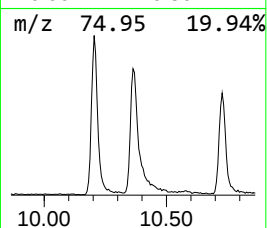
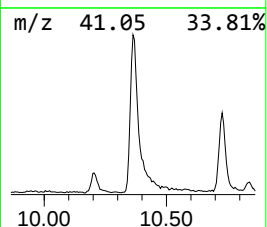
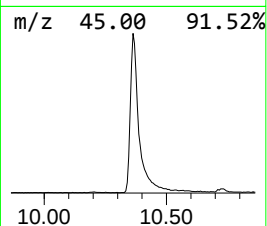
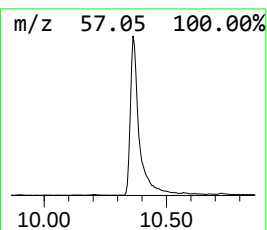
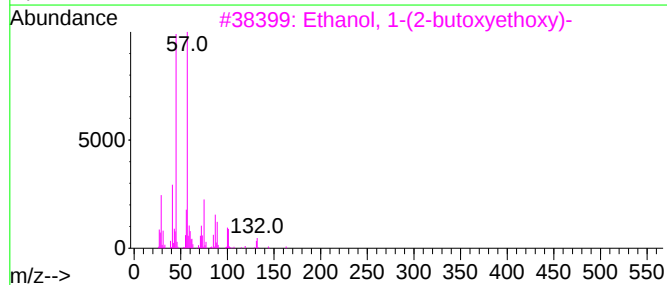
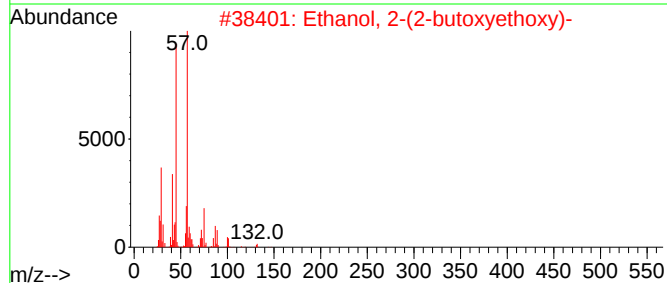
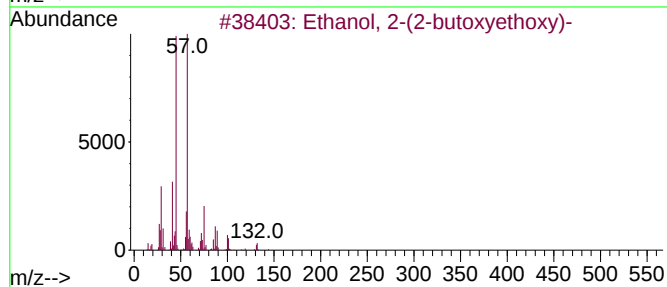
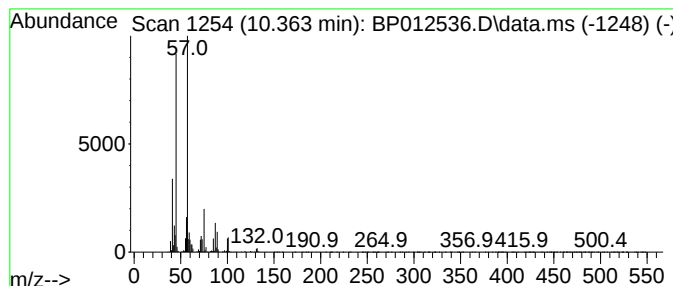
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.363	9.75 ng/ul	1178820	Naphthalene-d8	10.575

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, 1-(2-butoxyethoxy)-	162	C8H18O3	054446-78-5	83
4			Ethanol, 1-(2-butoxyethoxy)-	162	C8H18O3	054446-78-5	83
5			Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	72



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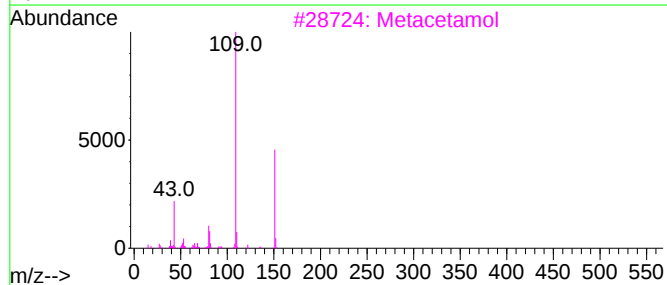
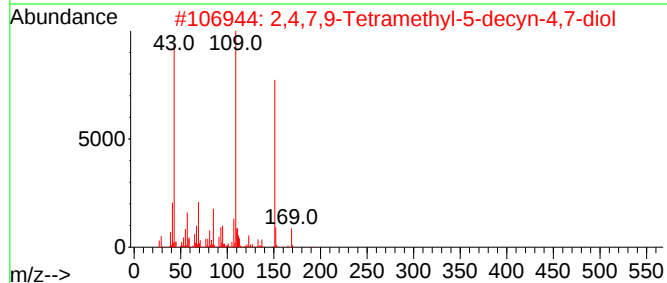
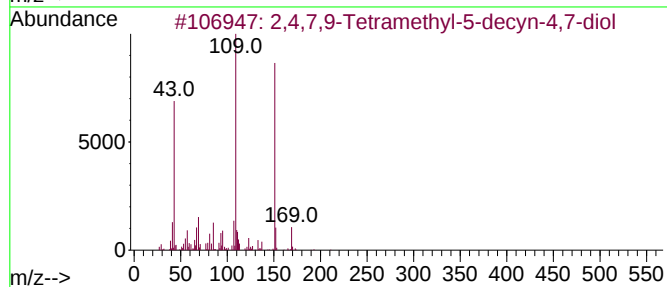
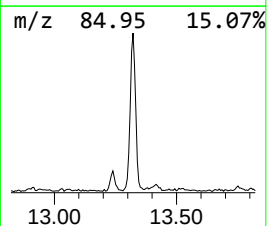
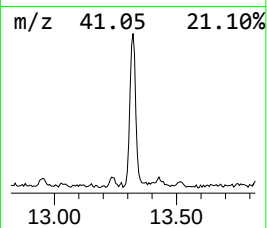
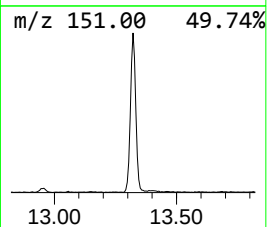
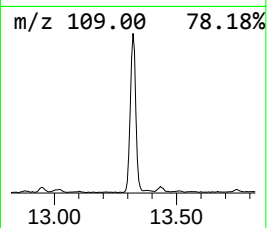
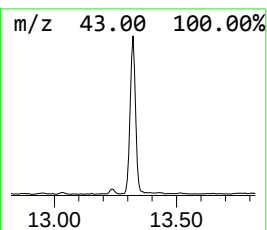
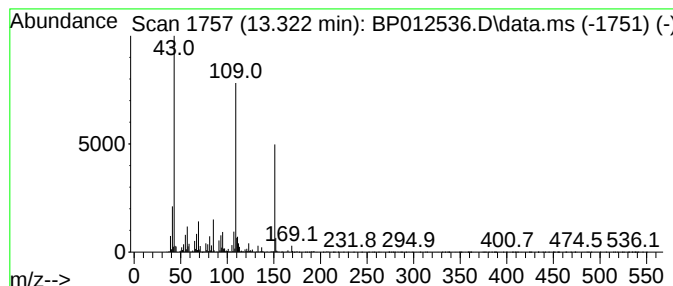
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.322	4.04 ng/ul	717409	Acenaphthene-d10	14.428

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	68		
3	Metacetamol	151	C8H9NO2	000621-42-1	64		
4	Acetaminophen	151	C8H9NO2	000103-90-2	59		
5	3-Pyridinol, 4-methyl-, acetate ...	151	C8H9NO2	001006-96-8	59		



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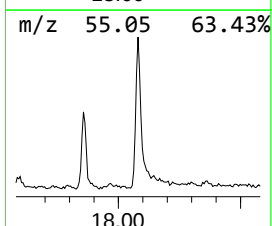
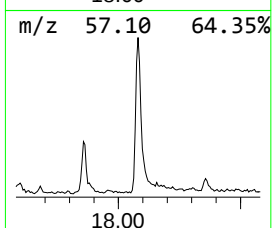
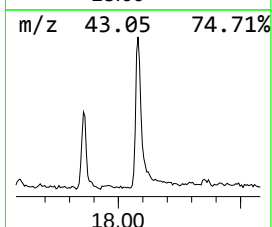
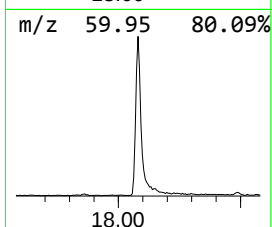
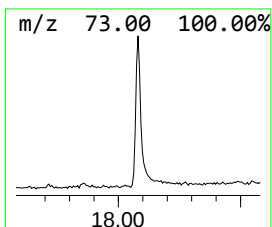
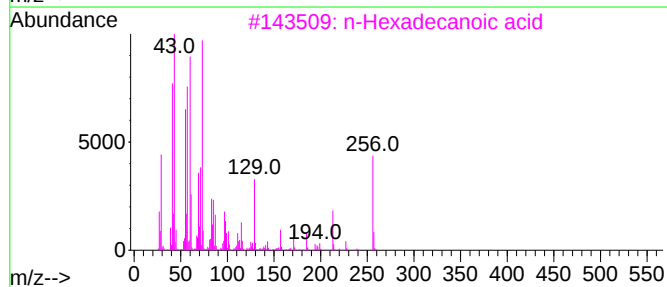
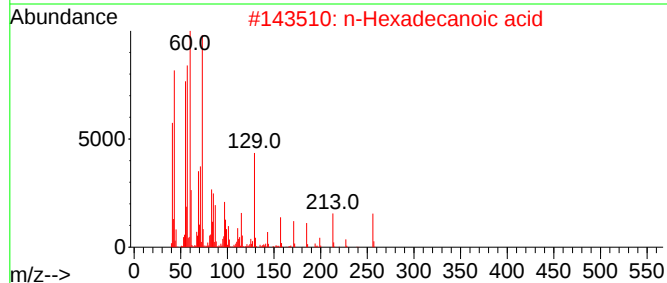
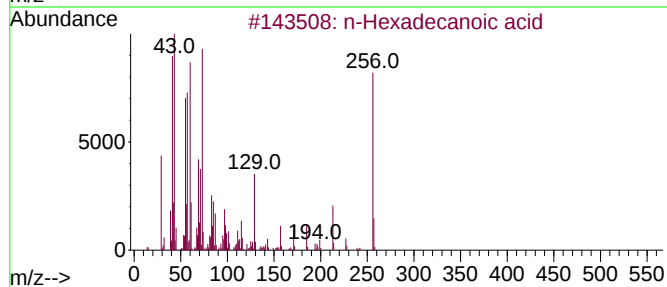
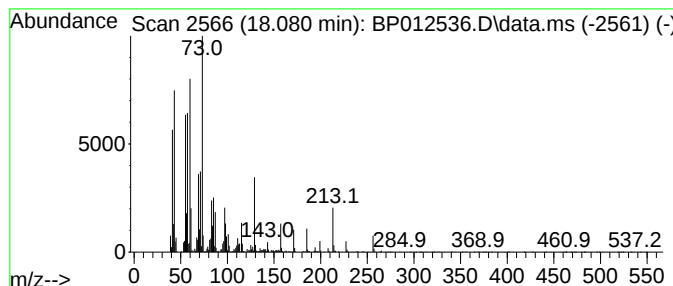
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Quant Title : SVOA CALIBRATION

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

Peak Number 3 n-Hexadecanoic acid Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.080	3.00 ng/ul	673190	Phenanthrene-d10	17.192

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	91
5			Pentadecanoic acid	242	C15H30O2	001002-84-2	91



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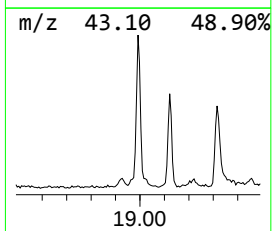
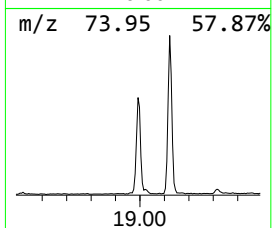
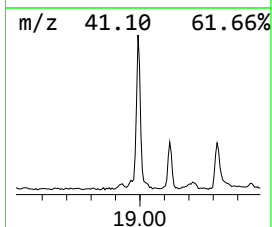
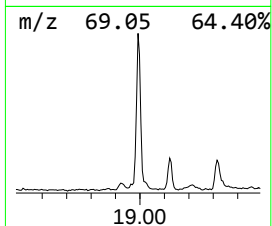
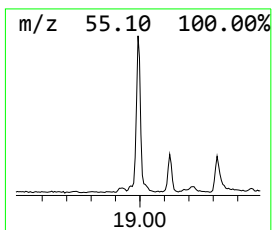
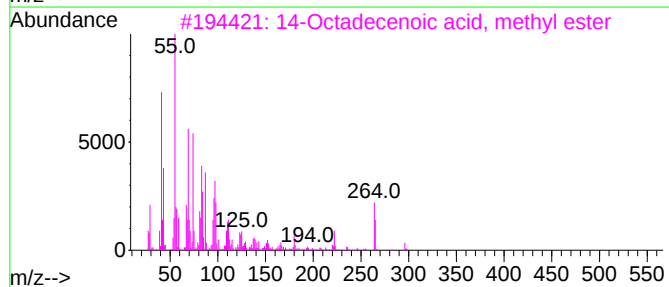
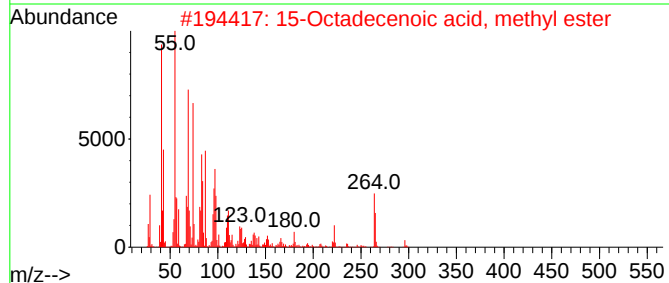
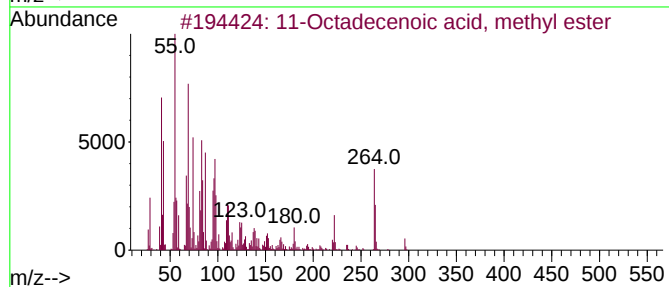
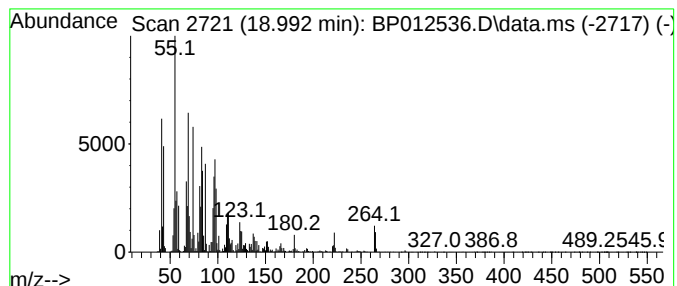
Instrument :
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ClientSampleId :
DBWX1

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
Quant Title : SVOA CALIBRATION

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

Peak Number 4 11-Octadecenoic acid, methyl... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.	
18.992	5.61 ng/ul	1257050	Phenanthrene-d10	17.192	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	11-Octadecenoic acid, methyl ester	296	C19H36O2	052380-33-3	99
2	15-Octadecenoic acid, methyl ester	296	C19H36O2	004764-72-1	99
3	14-Octadecenoic acid, methyl ester	296	C19H36O2	056554-48-4	99
4	6-Octadecenoic acid, methyl ester	296	C19H36O2	052355-31-4	99
5	16-Octadecenoic acid, methyl ester	296	C19H36O2	056554-49-5	99



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110722\
Data File : BP012536.D
Acq On : 07 Nov 2022 17:57
Operator : CG/JU
Sample : N5353-16
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DBWX1

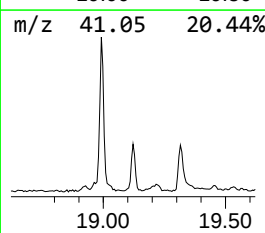
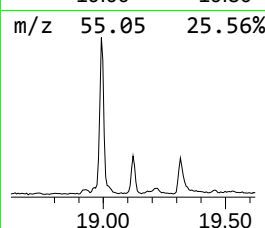
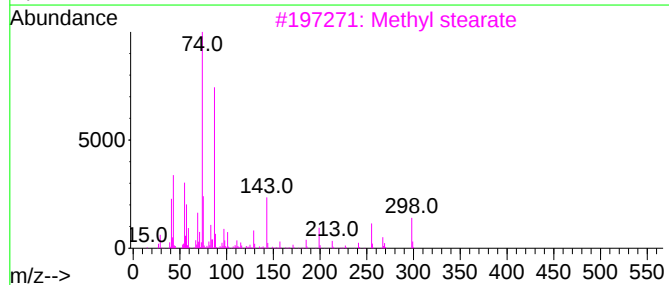
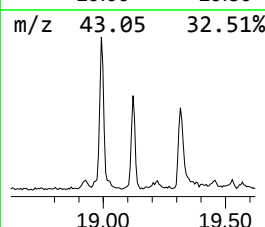
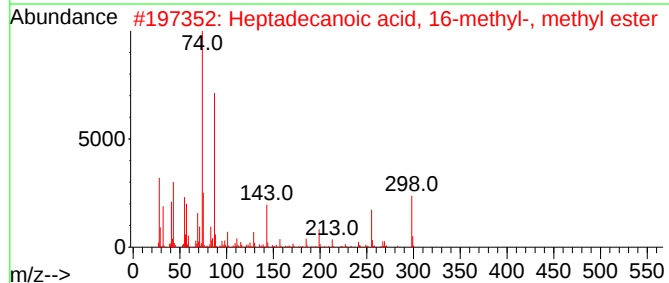
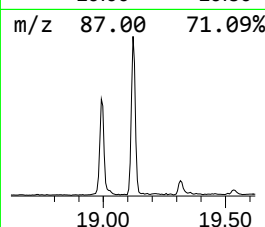
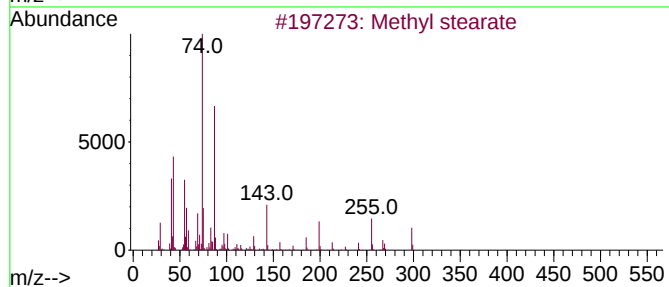
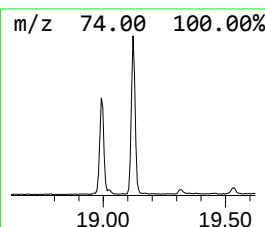
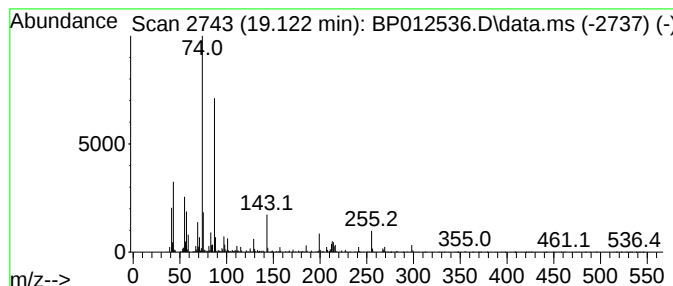
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 Methyl stearate Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.122	2.36 ng/ul	529832	Phenanthrene-d10	17.192

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Methyl stearate	298	C19H38O2	000112-61-8	97
2			Heptadecanoic acid, 16-methyl-, ...	298	C19H38O2	005129-61-3	92
3			Methyl stearate	298	C19H38O2	000112-61-8	92
4			Heptadecanoic acid, 15-methyl-, ...	298	C19H38O2	054833-55-5	91
5			Hexadecanoic acid, methyl ester	270	C17H34O2	000112-39-0	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110722\
Data File : BP012536.D
Acq On : 07 Nov 2022 17:57
Operator : CG/JU
Sample : N5353-16
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DBWX1

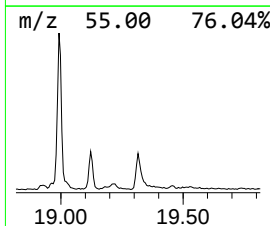
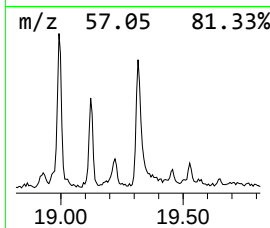
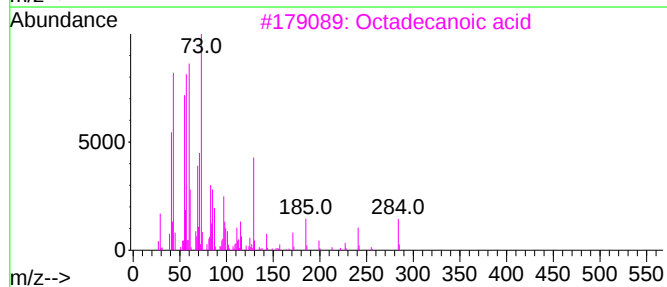
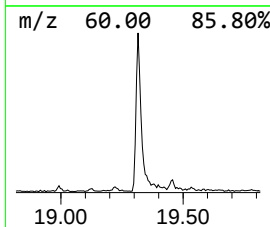
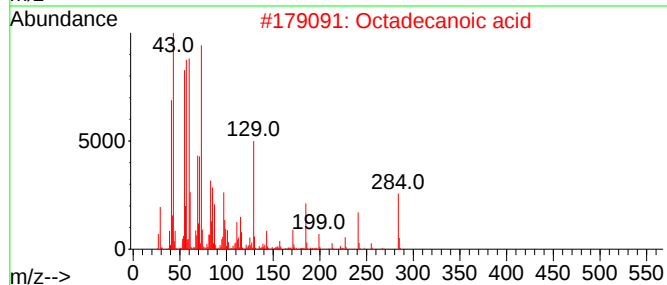
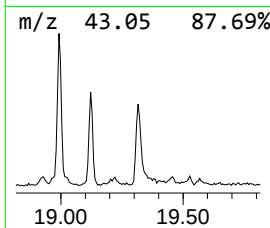
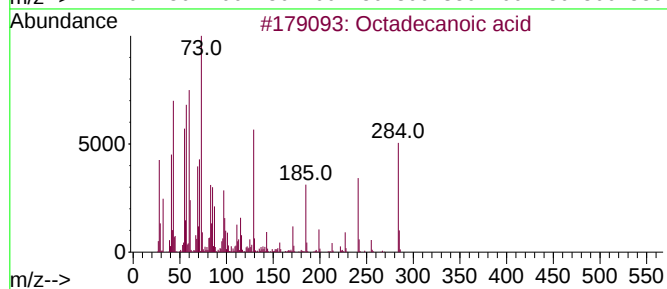
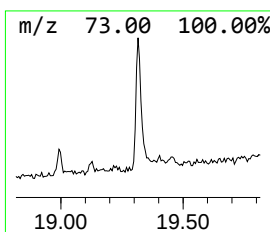
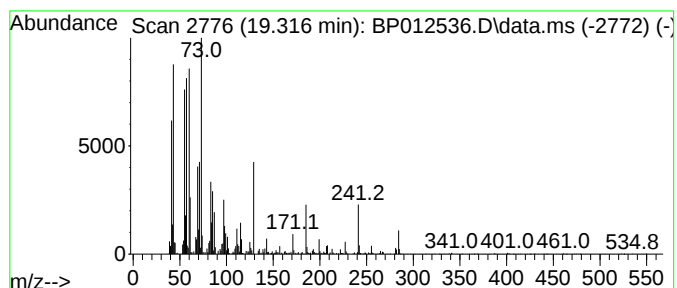
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
Quant Title : SVOA CALIBRATION

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

Peak Number 6 Octadecanoic acid Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.316	2.21 ng/ul	506224	Chrysene-d12	21.286

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	93
5			Pentadecanoic acid	242	C15H30O2	001002-84-2	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110722\
Data File : BP012536.D
Acq On : 07 Nov 2022 17:57
Operator : CG/JU
Sample : N5353-16
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DBWX1

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.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...	10.363	9.8	ng/ul	1178820	2	10.575	2417800	20.0
2,4,7,9-Tetrame...	13.322	4.0	ng/ul	717409	3	14.428	3548220	20.0
n-Hexadecanoic ...	18.080	3.0	ng/ul	673190	4	17.192	4481900	20.0
11-Octadecenoic...	18.992	5.6	ng/ul	1257050	4	17.192	4481900	20.0
Methyl stearate	19.122	2.4	ng/ul	529832	4	17.192	4481900	20.0
Octadecanoic acid	19.316	2.2	ng/ul	506224	5	21.286	4586360	20.0