

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041922\
 Data File : BP009965.D
 Acq On : 20 Apr 2022 02:29
 Operator : CG/JU
 Sample : N2422-12
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0J42

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-BP040722.M
 Title : SVOA CALIBRATION

Signal : TIC: BP009965.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.370	44	48	56	rVB	25838	39765	1.19%	0.105%
2	3.646	93	95	103	rVB4	10492	17168	0.51%	0.046%
3	3.793	117	120	137	rBV	110825	203395	6.09%	0.540%
4	4.123	173	176	181	rVB4	7256	9792	0.29%	0.026%
5	5.252	364	368	375	rBV	56096	85978	2.57%	0.228%
6	7.105	674	683	697	rVB2	96288	220371	6.59%	0.585%
7	7.269	704	711	720	rBV	310544	500573	14.98%	1.328%
8	7.340	720	723	730	rVB	14889	26681	0.80%	0.071%
9	7.475	739	746	757	rBV	301411	498966	14.93%	1.324%
10	7.946	817	826	835	rBV	432980	724268	21.67%	1.921%
11	8.011	835	837	841	rVV2	11457	16654	0.50%	0.044%
12	8.299	880	886	891	rBV3	25231	43158	1.29%	0.114%
13	8.646	936	945	954	rBV	271067	447894	13.40%	1.188%
14	9.099	1014	1022	1033	rBV	394355	666540	19.95%	1.768%
15	9.222	1035	1043	1047	rVV8	8505	20393	0.61%	0.054%
16	9.281	1047	1053	1057	rVV8	9907	21167	0.63%	0.056%
17	9.558	1095	1100	1106	rVB	29119	48289	1.45%	0.128%
18	9.828	1138	1146	1160	rBV	292907	526477	15.75%	1.397%
19	10.063	1180	1186	1189	rBV7	7303	12871	0.39%	0.034%
20	10.369	1229	1238	1255	rBV	448974	802640	24.02%	2.129%
21	10.528	1260	1265	1273	rVV	41848	78422	2.35%	0.208%
22	10.610	1275	1279	1285	rVB9	7675	14072	0.42%	0.037%
23	10.752	1295	1303	1311	rBV	641220	1083881	32.43%	2.875%
24	10.881	1317	1325	1343	rBV	637440	1147795	34.35%	3.045%
25	11.116	1360	1365	1367	rBV5	5716	10031	0.30%	0.027%
26	11.540	1432	1437	1446	rBV5	12358	24146	0.72%	0.064%
27	11.787	1472	1479	1483	rBV10	5272	11542	0.35%	0.031%
28	12.052	1520	1524	1531	rVB4	8108	11328	0.34%	0.030%
29	12.181	1542	1546	1551	rVB3	12169	18331	0.55%	0.049%
30	12.334	1565	1572	1580	rVB6	10596	24666	0.74%	0.065%
31	12.946	1671	1676	1679	rBV4	8378	13373	0.40%	0.035%
32	13.099	1699	1702	1705	rVV4	8638	11898	0.36%	0.032%
33	13.193	1712	1718	1724	rVB2	16757	25679	0.77%	0.068%
34	13.410	1751	1755	1760	rVV2	27207	43337	1.30%	0.115%
35	13.481	1762	1767	1774	rVV3	36627	63062	1.89%	0.167%
36	13.981	1845	1852	1858	rBV	1080542	1519582	45.47%	4.031%

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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 >

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Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP040722.M
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37	14.028	1858	1860	1865	rVV6	13588	22649	0.68%	0.060%
38	14.110	1871	1874	1883	rVB10	14901	25644	0.77%	0.068%
39	14.222	1888	1893	1895	rBV2	72781	98210	2.94%	0.261%
40	14.269	1895	1901	1915	rVB	1062270	1614176	48.30%	4.282%
41	14.481	1932	1937	1942	rVB	22147	33441	1.00%	0.089%
42	14.575	1946	1953	1964	rBV	1073889	1598255	47.83%	4.240%
43	14.716	1973	1977	1980	rBV6	7817	12472	0.37%	0.033%
44	14.757	1980	1984	1992	rBV2	82577	145045	4.34%	0.385%
45	14.993	2019	2024	2032	rVB2	9472	17153	0.51%	0.046%
46	15.210	2056	2061	2068	rVV2	46062	73186	2.19%	0.194%
47	15.381	2085	2090	2095	rVV2	72840	125249	3.75%	0.332%
48	15.434	2095	2099	2103	rVV	25288	36329	1.09%	0.096%
49	15.569	2115	2122	2135	rBV	1497599	2223328	66.53%	5.898%
50	15.675	2135	2140	2148	rBV	432892	612332	18.32%	1.624%
51	15.810	2158	2163	2166	rBV5	18369	26419	0.79%	0.070%
52	15.934	2177	2184	2188	rVB3	28605	45984	1.38%	0.122%
53	15.987	2188	2193	2200	rVB2	21910	32797	0.98%	0.087%
54	16.063	2201	2206	2211	rBV5	20176	31104	0.93%	0.083%
55	16.116	2211	2215	2223	rBV3	29721	50059	1.50%	0.133%
56	16.210	2227	2231	2235	rVV5	24235	36032	1.08%	0.096%
57	16.257	2235	2239	2243	rVV3	58699	99602	2.98%	0.264%
58	16.293	2243	2245	2248	rVV3	35489	47728	1.43%	0.127%
59	16.322	2248	2250	2253	rVV3	19989	27275	0.82%	0.072%
60	16.357	2254	2256	2262	rVB6	19304	28609	0.86%	0.076%
61	16.469	2270	2275	2279	rBV3	43193	69474	2.08%	0.184%
62	16.575	2288	2293	2297	rBV4	15273	23796	0.71%	0.063%
63	16.704	2312	2315	2316	rBV3	11473	10124	0.30%	0.027%
64	16.804	2328	2332	2336	rBV3	33411	49255	1.47%	0.131%
65	16.869	2341	2343	2347	rVB5	12754	14670	0.44%	0.039%
66	17.045	2370	2373	2375	rBV4	15204	18118	0.54%	0.048%
67	17.092	2375	2381	2383	rBV5	36592	69078	2.07%	0.183%
68	17.251	2404	2408	2414	rVB	107231	131675	3.94%	0.349%
69	17.328	2414	2421	2431	rBV	1322569	1982598	59.33%	5.259%
70	17.428	2431	2438	2445	rBV	1720163	2611805	78.16%	6.928%
71	17.616	2465	2470	2476	rBV	128924	204646	6.12%	0.543%
72	17.710	2482	2486	2493	rBV	96212	151239	4.53%	0.401%
73	17.945	2522	2526	2529	rBV5	16675	27682	0.83%	0.073%
74	17.998	2530	2535	2540	rVB	311492	378899	11.34%	1.005%
75	18.169	2562	2564	2567	rBV4	14439	18795	0.56%	0.050%
76	18.210	2567	2571	2577	rVV	240936	325620	9.74%	0.864%

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

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
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77	18.281	2578	2583	2587	rVV	304253	380443	11.38%	1.009%
78	18.328	2587	2591	2594	rVB4	11635	18118	0.54%	0.048%
79	18.628	2638	2642	2645	rBV	55132	68608	2.05%	0.182%
80	18.792	2667	2670	2674	rBV2	65490	77058	2.31%	0.204%
81	18.951	2694	2697	2702	rBV	67199	86769	2.60%	0.230%
82	19.039	2708	2712	2719	rBV	251729	343789	10.29%	0.912%
83	19.128	2719	2727	2730	rBV2	568837	759700	22.73%	2.015%
84	19.257	2745	2749	2754	rVB2	195320	250031	7.48%	0.663%
85	19.322	2757	2760	2761	rVV2	52146	57193	1.71%	0.152%
86	19.345	2761	2764	2770	rVB	222632	272282	8.15%	0.722%
87	19.439	2776	2780	2783	rBV3	55535	69919	2.09%	0.185%
88	19.569	2799	2802	2808	rVB2	75895	102631	3.07%	0.272%
89	19.651	2810	2816	2826	rBV	2397760	3341791	100.00%	8.865%
90	20.551	2967	2969	2975	rVB	50400	64565	1.93%	0.171%
91	20.757	3001	3004	3009	rBV2	111303	152181	4.55%	0.404%
92	20.804	3009	3012	3016	rVV	111788	137059	4.10%	0.364%
93	20.863	3019	3022	3029	rVB2	59874	78568	2.35%	0.208%
94	21.110	3060	3064	3073	rVB	1040654	1199354	35.89%	3.182%
95	21.239	3083	3086	3089	rVB	77758	78760	2.36%	0.209%
96	21.316	3093	3099	3103	rBV3	100423	168037	5.03%	0.446%
97	21.404	3109	3114	3129	rVB	1538460	2187239	65.45%	5.802%
98	23.704	3497	3505	3512	rBV2	1530356	3117597	93.29%	8.270%
99	23.857	3524	3531	3544	rBV2	1023676	2189926	65.53%	5.809%
100	24.604	3653	3658	3668	rVB3	136408	310731	9.30%	0.824%

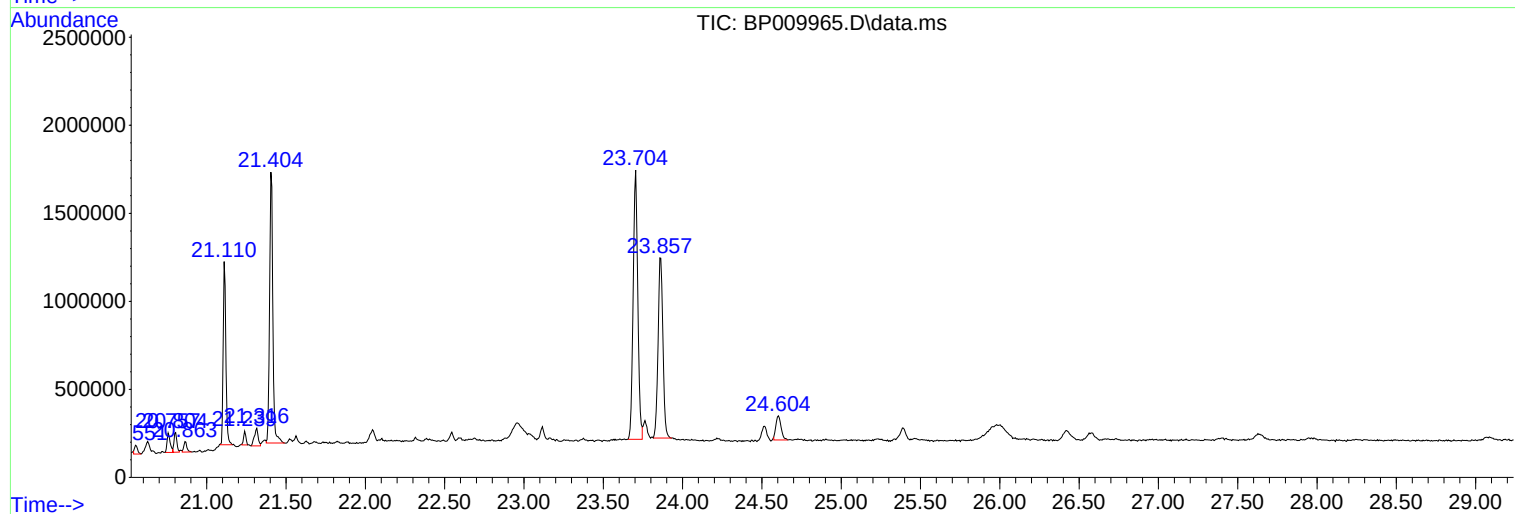
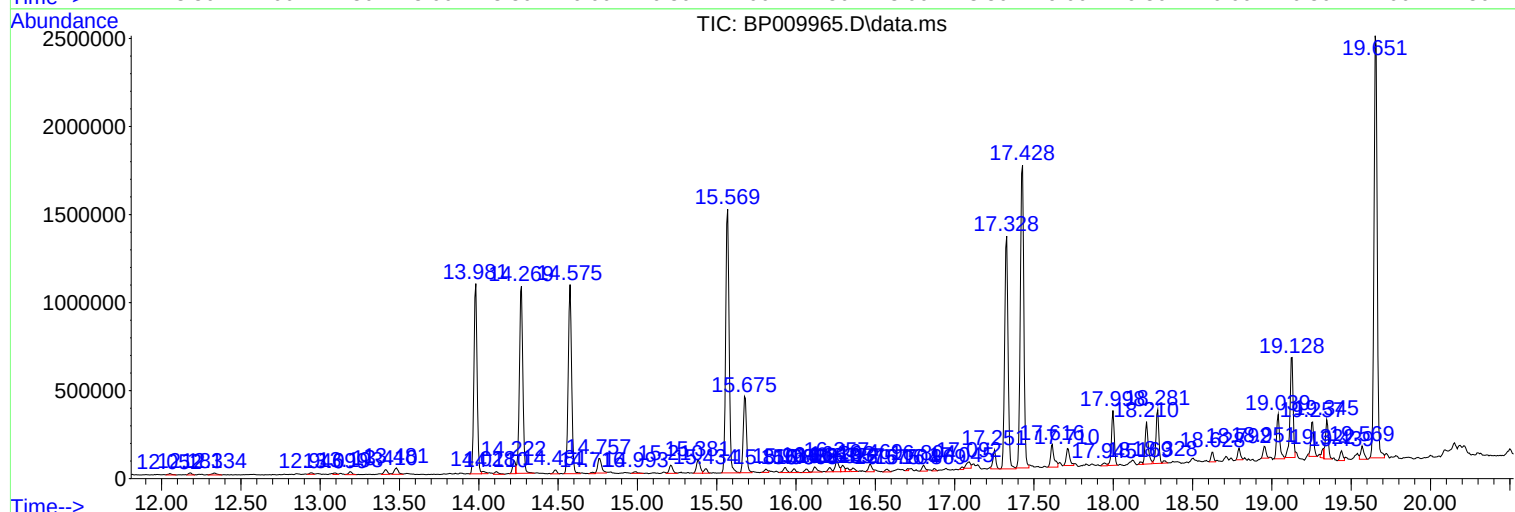
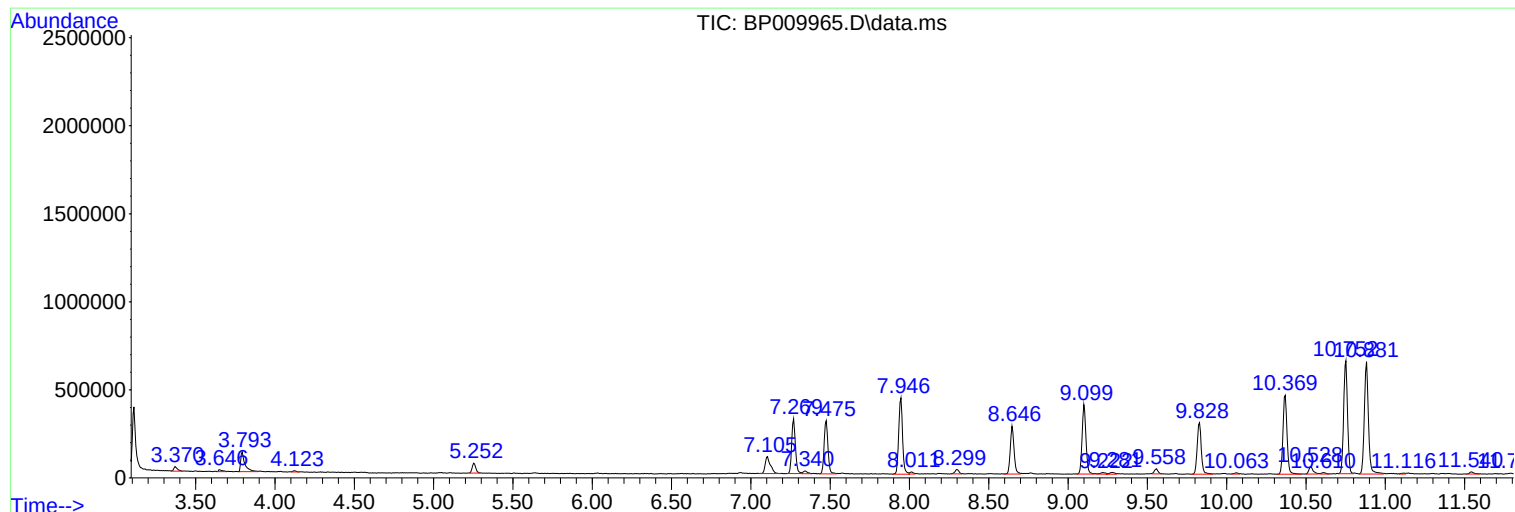
Sum of corrected areas: 37697086

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP040722.M
Quant Title : SVOA CALIBRATION

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



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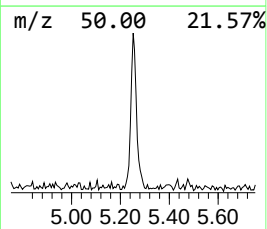
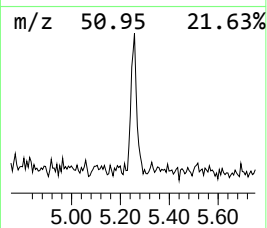
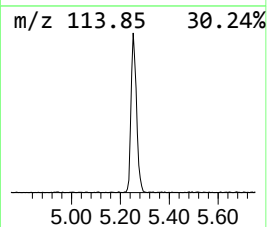
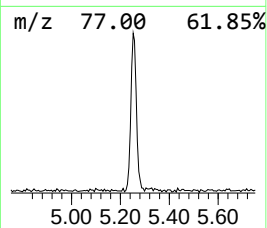
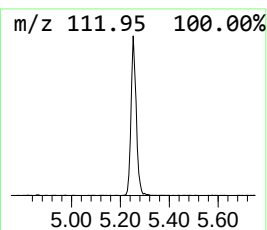
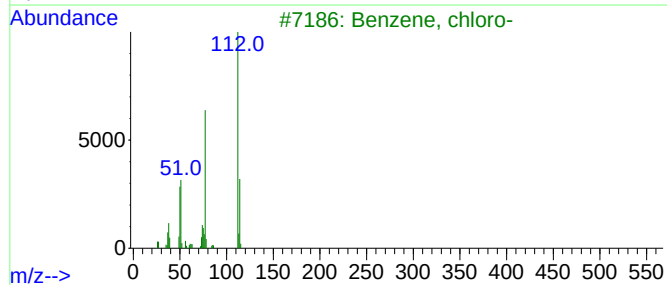
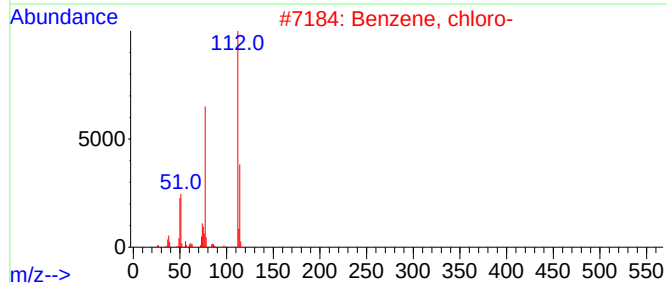
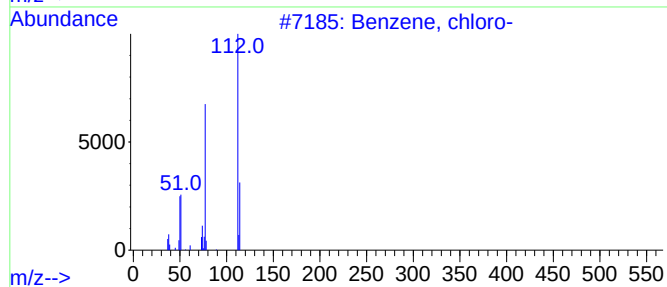
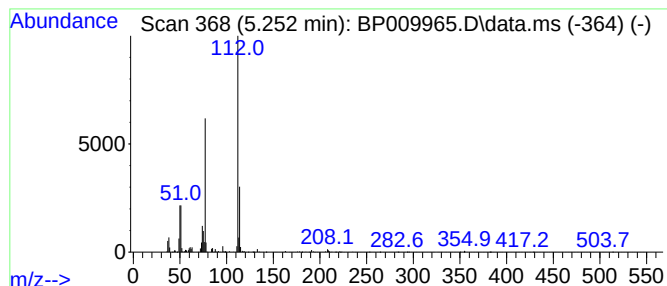
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP040722.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Benzene, chloro- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.252	2.37 ng/ul	85978	1,4-Dichlorobenzene-d4	7.946

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, chloro-	112	C6H5Cl	000108-90-7	96
2			Benzene, chloro-	112	C6H5Cl	000108-90-7	95
3			Benzene, chloro-	112	C6H5Cl	000108-90-7	91
4			Benzene, chloro-	112	C6H5Cl	000108-90-7	91
5			Benzene, chloro-	112	C6H5Cl	000108-90-7	91



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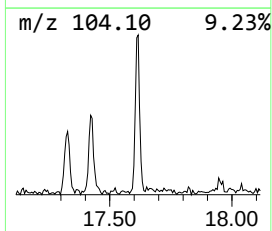
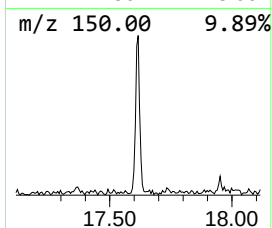
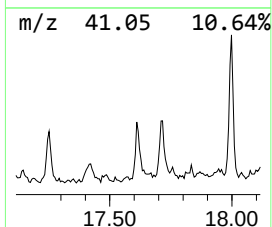
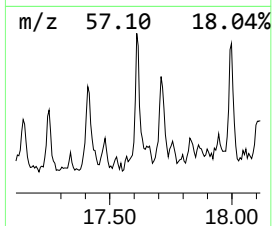
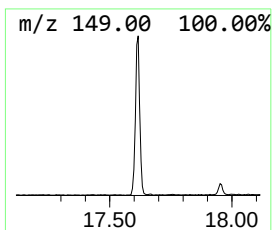
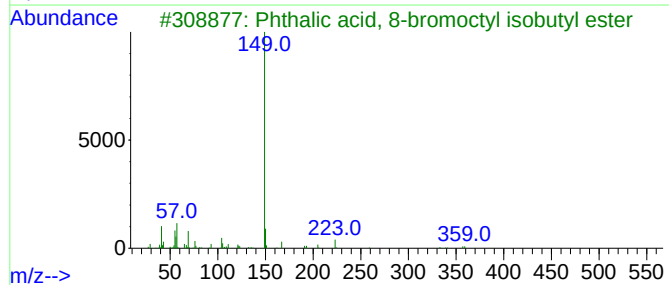
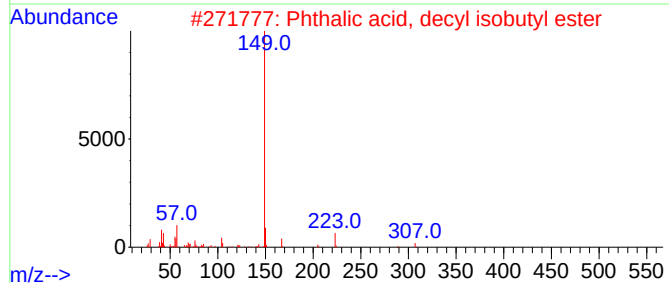
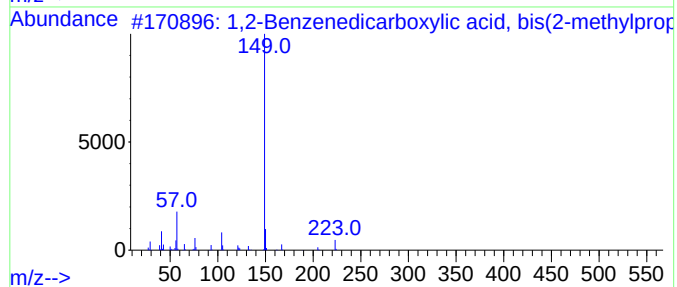
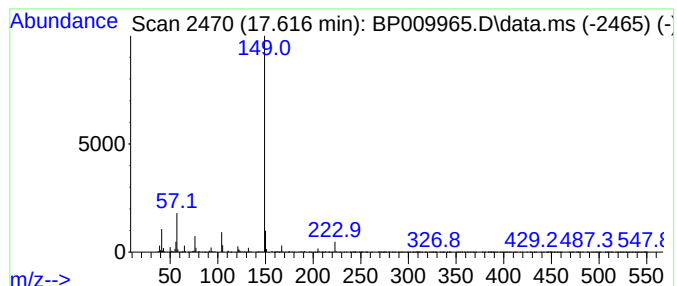
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP040722.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 1,2-Benzenedicarboxylic aci... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.616	2.06 ng/ul	204646	Phenanthrene-d10	17.328

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1,2-Benzenedicarboxylic acid, bi...	278	C16H22O4	000084-69-5	91
2			Phthalic acid, decyl isobutyl ester	362	C22H34O4	1000308-94-2	83
3			Phthalic acid, 8-bromooctyl isobu...	412	C20H29BrO4	1000382-55-3	83
4			Phthalic acid, butyl isohexyl ester	306	C18H26O4	1000309-03-6	78
5			Phthalic acid, isobutyl 2-methyl...	312	C19H20O4	1000314-95-7	72



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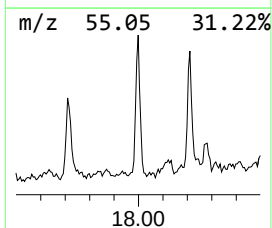
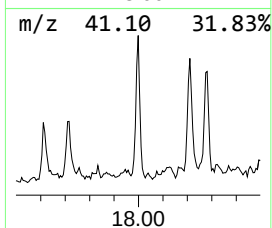
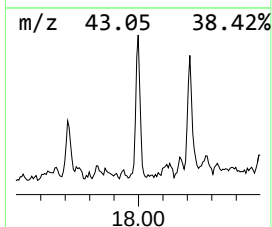
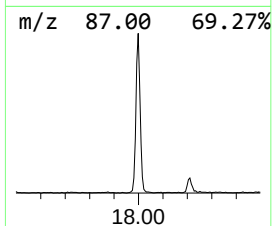
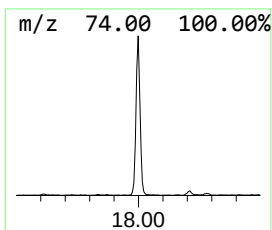
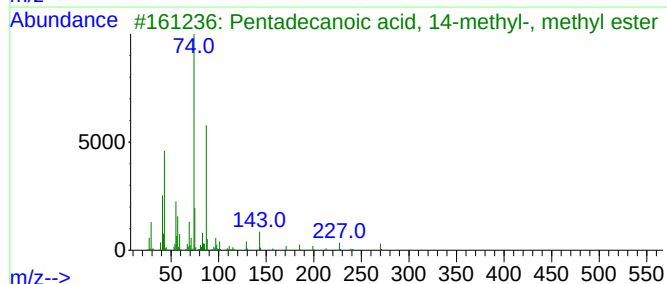
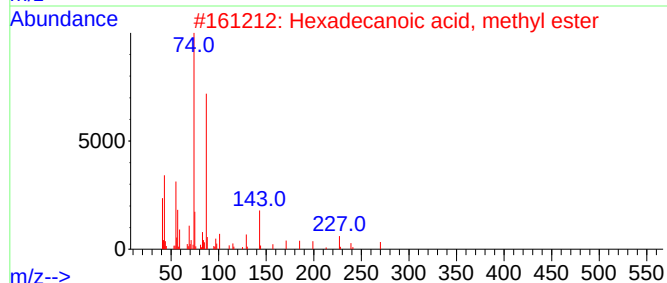
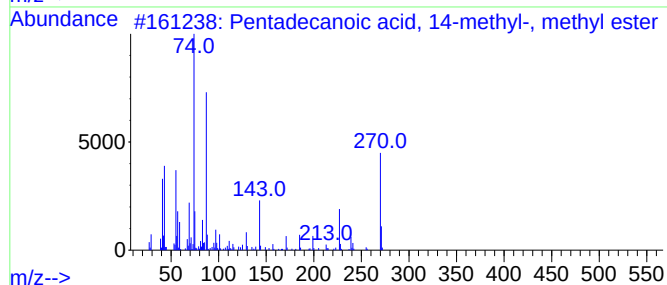
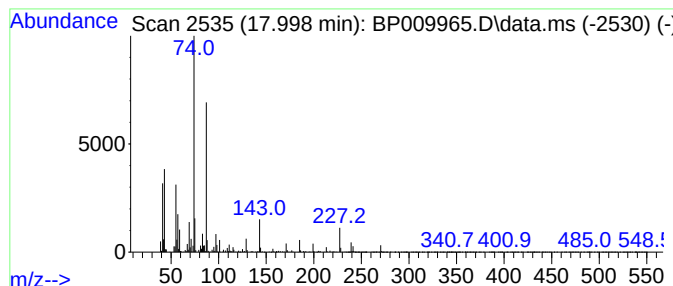
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Quant Title : SVOA CALIBRATION

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

Peak Number 3 Pentadecanoic acid, 14-meth... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.998	3.82 ng/ul	378899	Phenanthrene-d10	17.328

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentadecanoic acid, 14-methyl-, ...	270	C17H34O2	005129-60-2	97
2			Hexadecanoic acid, methyl ester	270	C17H34O2	000112-39-0	96
3			Pentadecanoic acid, 14-methyl-, ...	270	C17H34O2	005129-60-2	95
4			Tridecanoic acid, methyl ester	228	C14H28O2	001731-88-0	93
5			Hexadecanoic acid, methyl ester	270	C17H34O2	000112-39-0	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041922\
Data File : BP009965.D
Acq On : 20 Apr 2022 02:29
Operator : CG/JU
Sample : N2422-12
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0J42

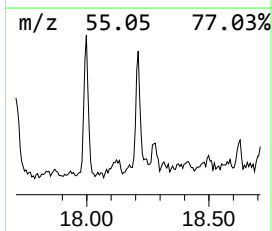
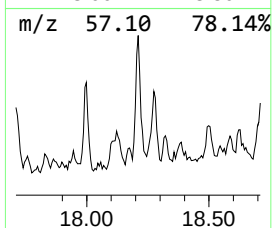
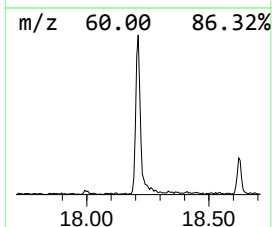
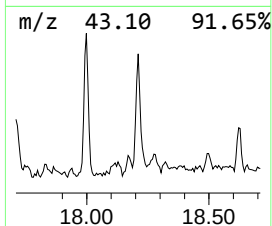
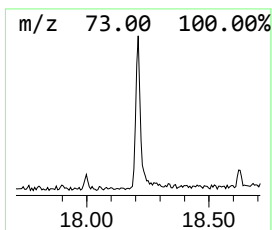
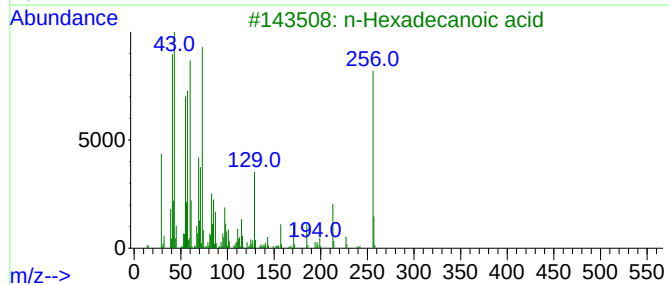
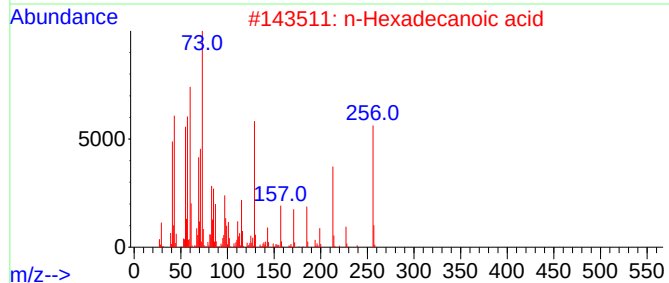
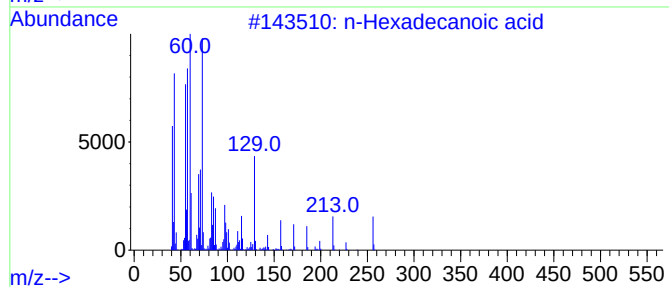
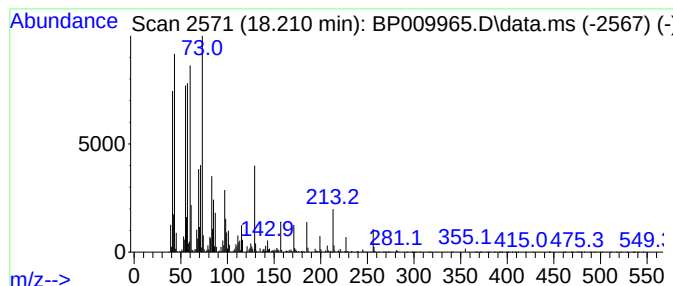
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 n-Hexadecanoic acid Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.210	3.28 ng/ul	325620	Phenanthrene-d10	17.328

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	98
4			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	97
5			Pentadecanoic acid	242	C15H30O2	001002-84-2	83



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041922\
Data File : BP009965.D
Acq On : 20 Apr 2022 02:29
Operator : CG/JU
Sample : N2422-12
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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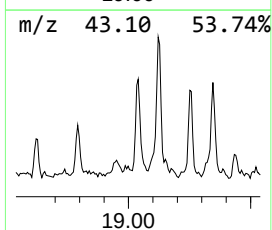
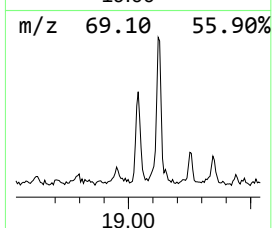
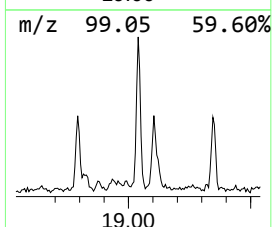
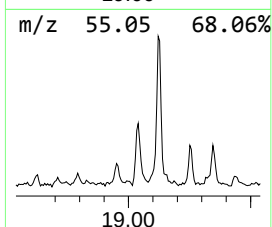
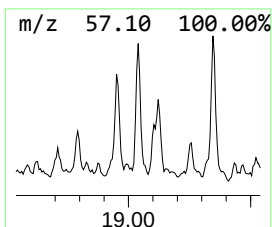
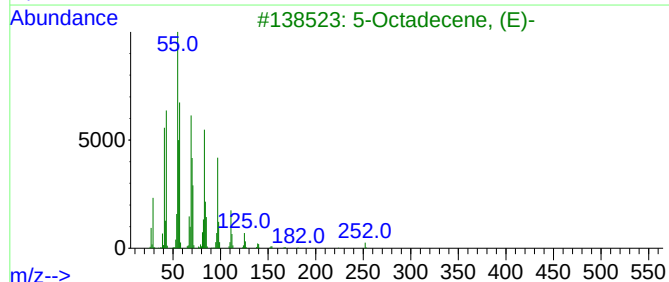
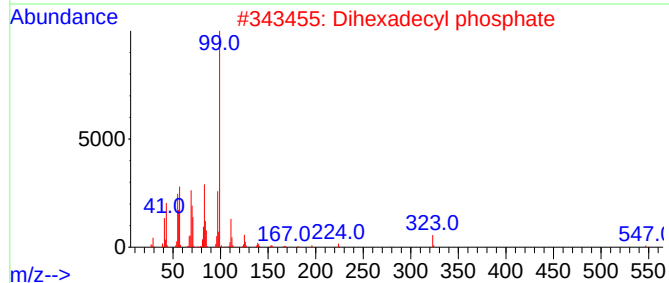
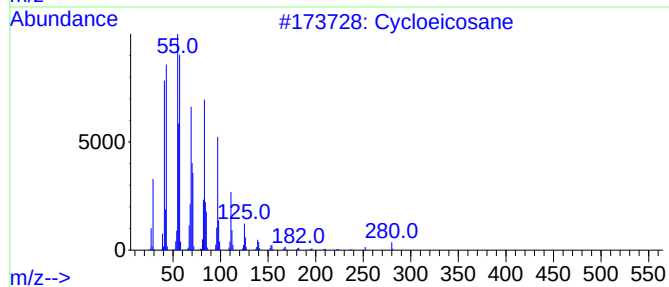
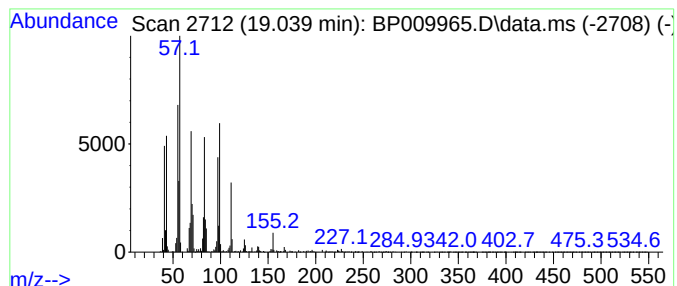
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 (DEL) Alkane: Cyclic19.040 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.040	3.47 ng/ul	343789	Phenanthrene-d10	17.328

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cycloeicosane	280	C20H40	000296-56-0	49
2			Dihexadecyl phosphate	546	C32H67O4P	002197-63-9	49
3			5-Octadecene, (E)-	252	C18H36	007206-21-5	46
4			E-15-Heptadecenal	252	C17H32O	1000130-97-9	46
5			Cyclopentane, 2-isopropyl-1,3-di...	140	C10H20	032281-85-9	46



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041922\
Data File : BP009965.D
Acq On : 20 Apr 2022 02:29
Operator : CG/JU
Sample : N2422-12
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0J42

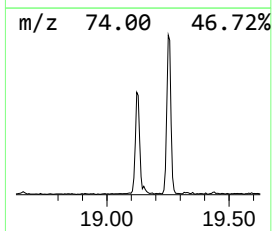
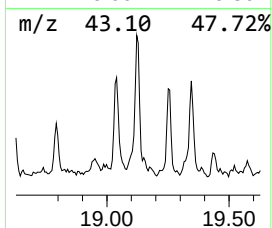
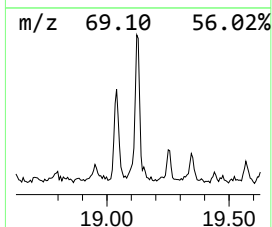
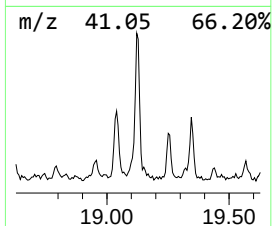
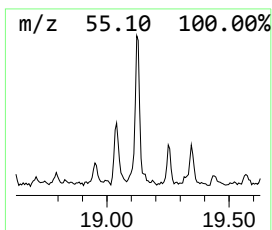
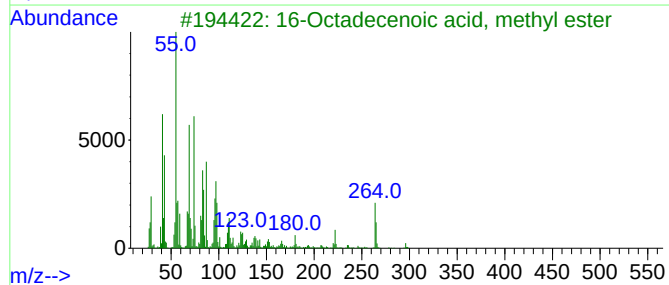
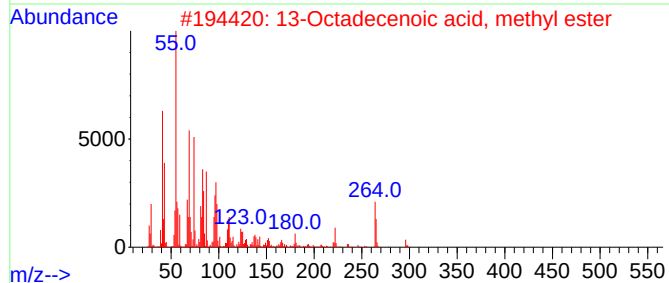
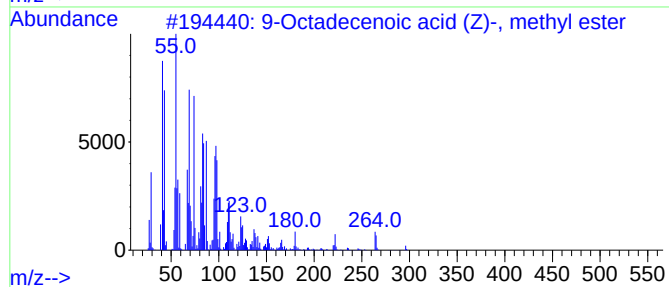
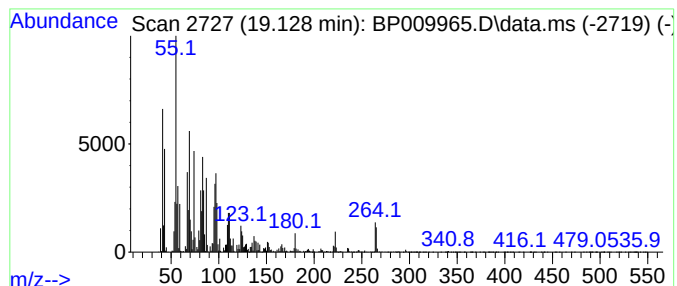
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 9-Octadecenoic acid (Z)-, m... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.128	7.66 ng/ul	759700	Phenanthrene-d10	17.328

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			9-Octadecenoic acid (Z)-, methyl...	296	C19H36O2	000112-62-9	98
2			13-Octadecenoic acid, methyl ester	296	C19H36O2	056554-47-3	97
3			16-Octadecenoic acid, methyl ester	296	C19H36O2	056554-49-5	95
4			14-Octadecenoic acid, methyl ester	296	C19H36O2	056554-48-4	95
5			cis-13-Octadecenoic acid, methyl...	296	C19H36O2	1010333-58-3	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041922\
Data File : BP009965.D
Acq On : 20 Apr 2022 02:29
Operator : CG/JU
Sample : N2422-12
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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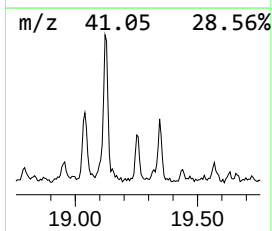
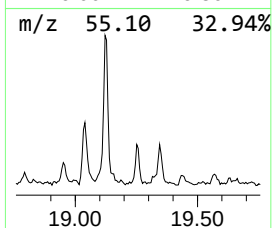
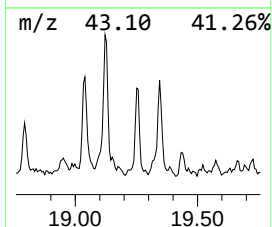
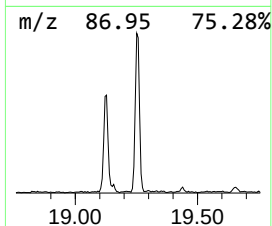
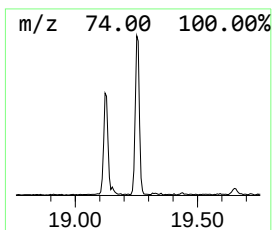
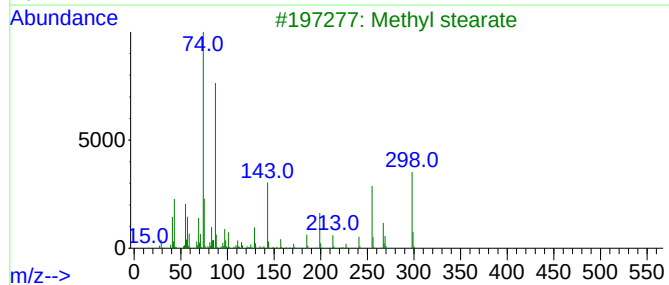
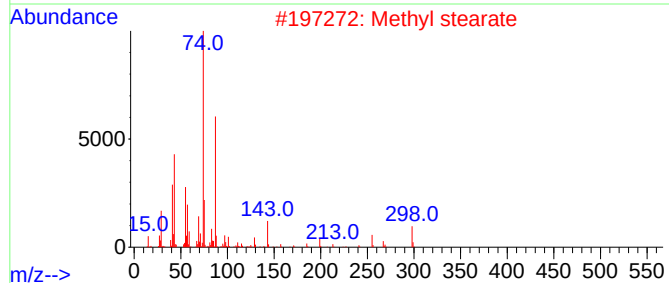
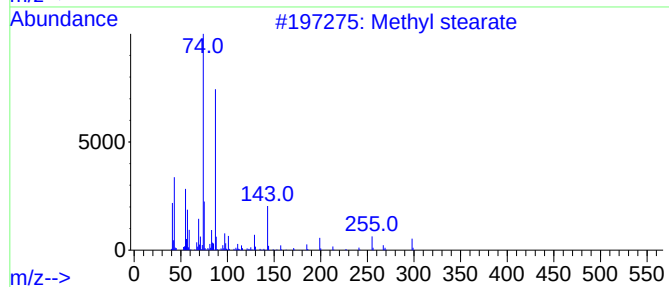
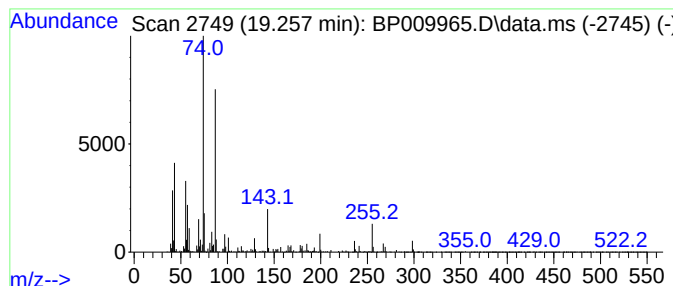
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 Methyl stearate Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.257	2.52 ng/ul	250031	Phenanthrene-d10	17.328

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Methyl stearate	298	C19H38O2	000112-61-8	98
2			Methyl stearate	298	C19H38O2	000112-61-8	97
3			Methyl stearate	298	C19H38O2	000112-61-8	95
4			Methyl stearate	298	C19H38O2	000112-61-8	95
5			Methyl stearate	298	C19H38O2	000112-61-8	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041922\
Data File : BP009965.D
Acq On : 20 Apr 2022 02:29
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Sample : N2422-12
Misc :
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Instrument :
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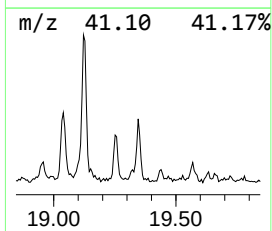
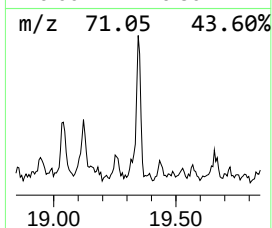
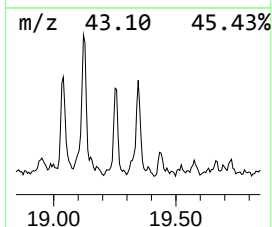
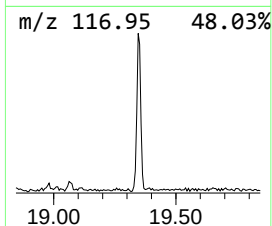
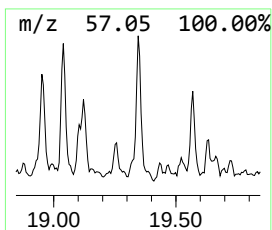
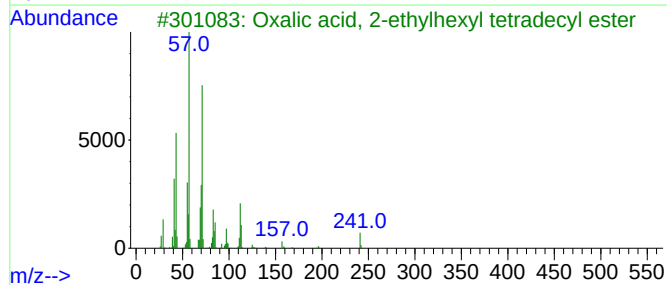
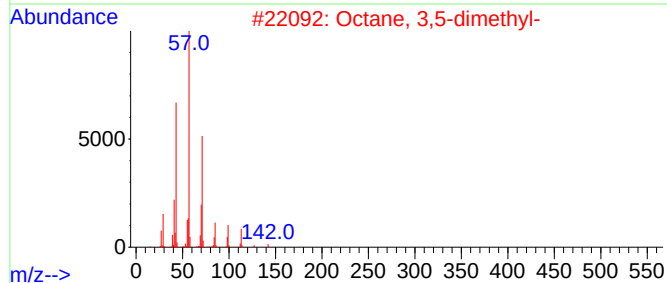
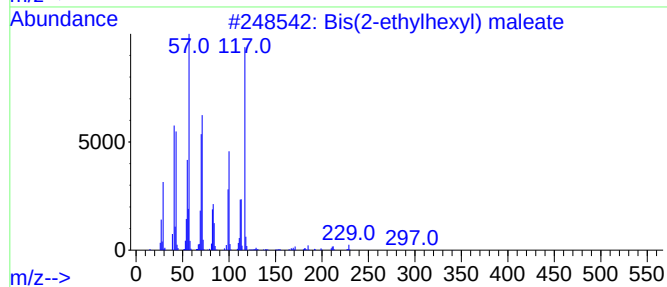
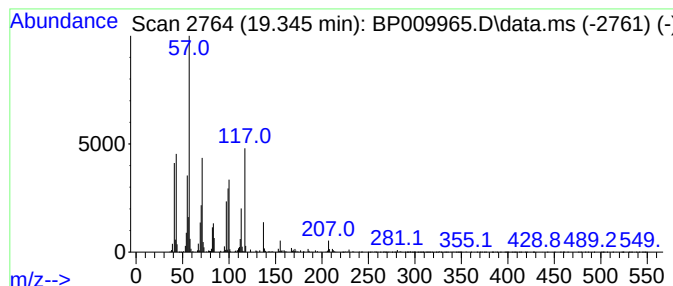
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 unknown-01 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.345	2.75 ng/ul	272282	Phenanthrene-d10	17.328

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Bis(2-ethylhexyl) maleate	340	C20H36O4	000142-16-5	43
2			Octane, 3,5-dimethyl-	142	C10H22	015869-93-9	27
3			Oxalic acid, 2-ethylhexyl tetrad...	398	C24H46O4	1000309-39-7	27
4			Sulfurous acid, butyl 2-ethylhex...	250	C12H26O3S	1000309-18-8	27
5			Pentadecane, 2,6,10-trimethyl-	254	C18H38	003892-00-0	25



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041922\
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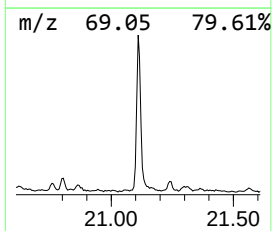
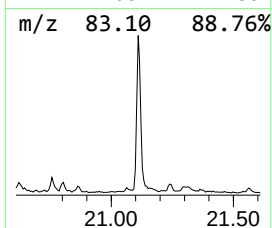
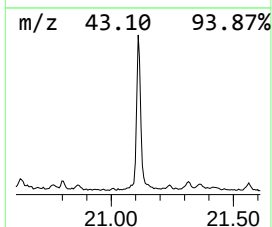
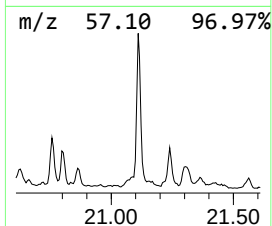
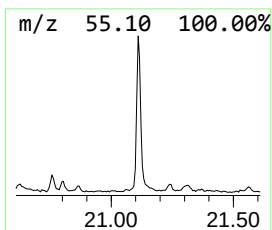
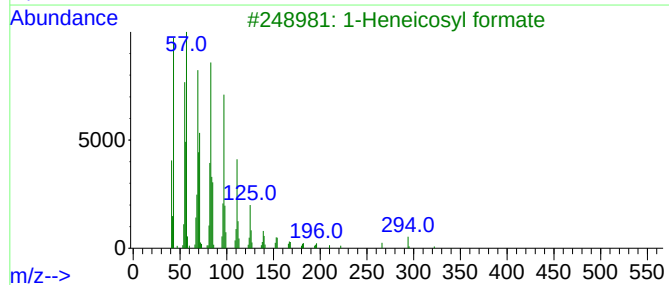
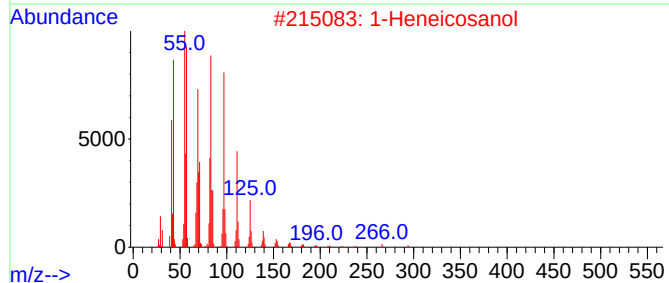
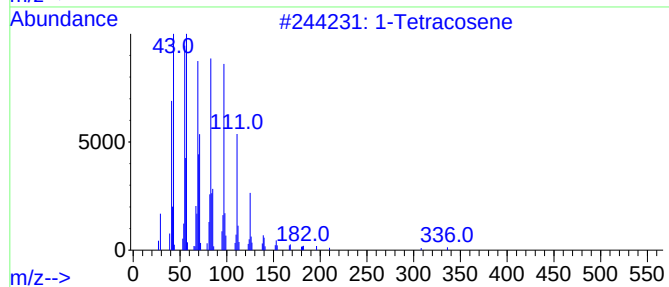
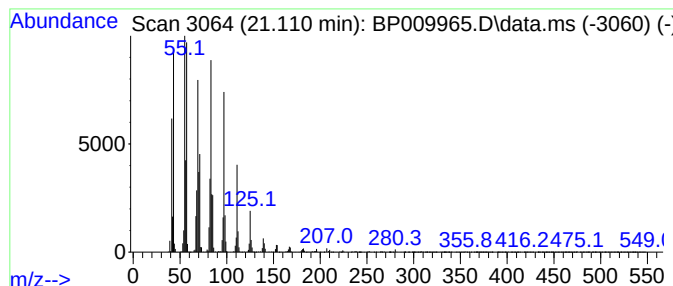
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 (DEL) Alkane: Straight-Chai... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.	
21.110	10.97 ng/ul	1199350	Chrysene-d12	21.404	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Tetracosene	336	C24H48	010192-32-2	95
2	1-Heneicosanol	312	C21H44O	015594-90-8	95
3	1-Heneicosyl formate	340	C22H44O2	077899-03-7	94
4	1-Hexacosene	364	C26H52	018835-33-1	94
5	Behenic alcohol	326	C22H46O	000661-19-8	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041922\
Data File : BP009965.D
Acq On : 20 Apr 2022 02:29
Operator : CG/JU
Sample : N2422-12
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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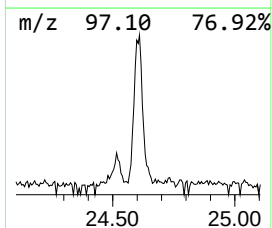
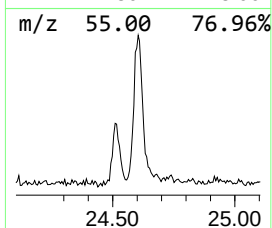
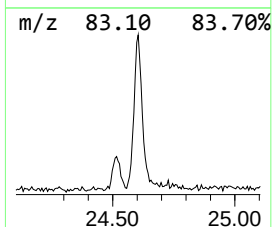
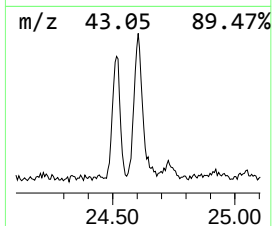
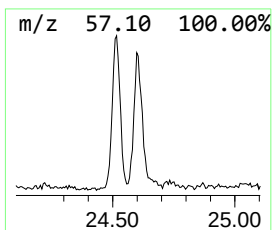
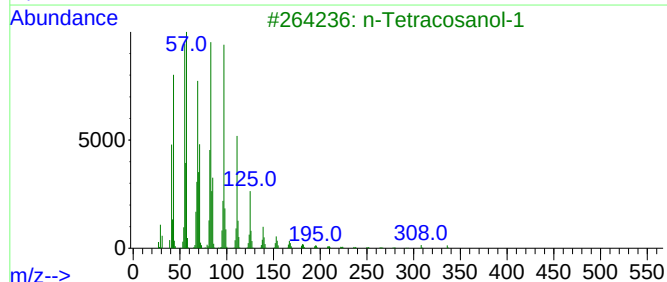
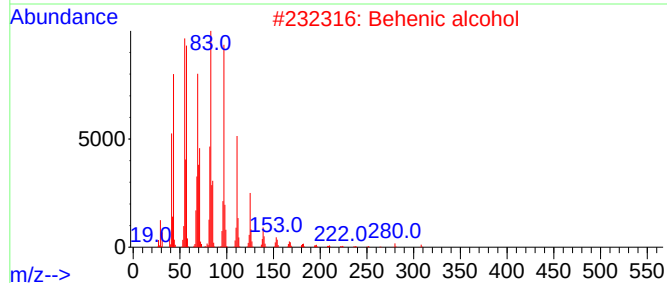
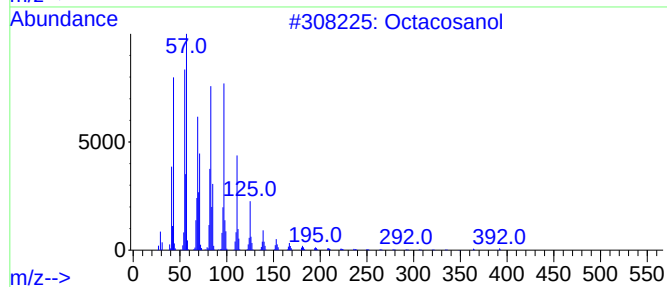
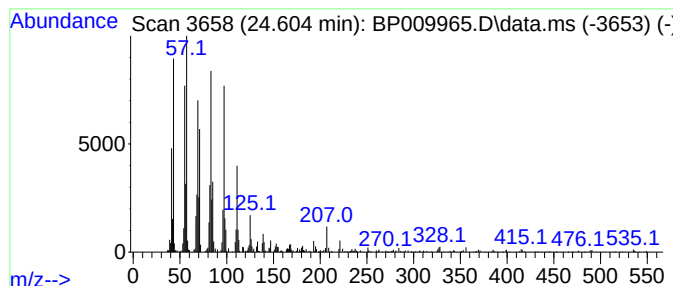
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP040722.M
Quant Title : SVOA CALIBRATION

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

Peak Number 10 Octacosanol Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.604	2.84 ng/ul	310731	Perylene-d12	23.857

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octacosanol	410	C28H58O	000557-61-9	93
2			Behenic alcohol	326	C22H46O	000661-19-8	93
3			n-Tetracosanol-1	354	C24H50O	000506-51-4	93
4			1-Docosanol, methyl ether	340	C23H48O	1000333-91-9	93
5			1-Heneicosanol	312	C21H44O	015594-90-8	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041922\
Data File : BP009965.D
Acq On : 20 Apr 2022 02:29
Operator : CG/JU
Sample : N2422-12
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0J42

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP040722.M
Quant Title : SVOA CALIBRATION

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Benzene, chloro-	5.252	2.4	ng/ul	85978	1	7.946	724268	20.0
1,2-Benzenedica...	17.616	2.1	ng/ul	204646	4	17.328	1982600	20.0
Pentadecanoic a...	17.998	3.8	ng/ul	378899	4	17.328	1982600	20.0
n-Hexadecanoic ...	18.210	3.3	ng/ul	325620	4	17.328	1982600	20.0
(DEL) Alkane: C...	19.040	3.5	ng/ul	343789	4	17.328	1982600	20.0
9-Octadecenoic ...	19.128	7.7	ng/ul	759700	4	17.328	1982600	20.0
Methyl stearate	19.257	2.5	ng/ul	250031	4	17.328	1982600	20.0
unknown-01	19.345	2.8	ng/ul	272282	4	17.328	1982600	20.0
(DEL) Alkane: S...	21.110	11.0	ng/ul	1199350	5	21.404	2187240	20.0
Octacosanol	24.604	2.8	ng/ul	310731	6	23.857	2189930	20.0