

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP112423\
 Data File : BP018333.D
 Acq On : 24 Nov 2023 18:26
 Operator : MA/JU
 Sample : 05264-09
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 DCKP7

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-BP112223.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BP018333.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.270	27	31	38	rVB	89618	113369	1.44%	0.113%
2	3.687	98	102	113	rBV	450906	622403	7.88%	0.621%
3	3.934	139	144	149	rVB2	22885	36467	0.46%	0.036%
4	4.028	157	160	164	rBV2	20791	27386	0.35%	0.027%
5	4.969	317	320	327	rVB8	7839	14178	0.18%	0.014%
6	5.575	421	423	427	rVB4	9366	9616	0.12%	0.010%
7	6.758	618	624	625	rBV6	5063	8011	0.10%	0.008%
8	7.028	663	670	678	rVB	375088	576053	7.29%	0.575%
9	7.205	691	700	708	rBV	1425572	2151076	27.24%	2.147%
10	7.399	725	733	743	rBV	1284244	1996081	25.27%	1.992%
11	7.499	745	750	757	rVB8	9229	17292	0.22%	0.017%
12	7.869	806	813	821	rBV	1368631	2094520	26.52%	2.090%
13	8.116	850	855	857	rBV6	4634	9298	0.12%	0.009%
14	8.575	924	933	945	rBV	1011161	1594322	20.19%	1.591%
15	9.028	1002	1010	1021	rBV	1514613	2420728	30.65%	2.416%
16	9.469	1076	1085	1093	rBV2	92264	171162	2.17%	0.171%
17	9.752	1124	1133	1144	rBV	867555	1385823	17.55%	1.383%
18	9.981	1164	1172	1181	rBV2	41443	88741	1.12%	0.089%
19	10.287	1215	1224	1235	rBV	1735726	2807739	35.55%	2.802%
20	10.669	1282	1289	1299	rVB	1728281	2870206	36.34%	2.864%
21	10.804	1302	1312	1321	rBV	1724139	2848129	36.06%	2.842%
22	11.169	1366	1374	1378	rBV5	7424	16923	0.21%	0.017%
23	11.451	1417	1422	1432	rBV4	22096	50700	0.64%	0.051%
24	11.963	1502	1509	1515	rVB2	66763	112625	1.43%	0.112%
25	12.257	1553	1559	1565	rBV	36955	58308	0.74%	0.058%
26	12.698	1627	1634	1638	rBV10	8970	21175	0.27%	0.021%
27	12.893	1660	1667	1670	rBV9	8403	20698	0.26%	0.021%
28	13.122	1705	1706	1712	rVB6	8751	11554	0.15%	0.012%
29	13.410	1751	1755	1760	rVB4	21769	35231	0.45%	0.035%
30	13.469	1760	1765	1772	rVB8	7954	15874	0.20%	0.016%
31	13.822	1821	1825	1832	rVB9	7473	18579	0.24%	0.019%
32	13.916	1835	1841	1854	rVV	3245911	4468109	56.57%	4.459%
33	14.193	1880	1888	1897	rBV	3781566	5521333	69.91%	5.510%
34	14.498	1932	1940	1947	rBV	2338353	3406988	43.14%	3.400%
35	14.687	1966	1972	1984	rBV	134795	223386	2.83%	0.223%
36	14.922	2006	2012	2018	rBV2	57468	93225	1.18%	0.093%

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37	15.157	2049	2052	2057	rVB7	7370	8652	0.11%	0.009%
38	15.316	2071	2079	2085	rBV2	33003	55529	0.70%	0.055%
39	15.492	2102	2109	2118	rBV	4619506	6722026	85.11%	6.709%
40	15.604	2122	2128	2139	rBV	147610	257980	3.27%	0.257%
41	15.728	2146	2149	2156	rVB4	21558	33494	0.42%	0.033%
42	15.916	2178	2181	2183	rBV4	7906	9992	0.13%	0.010%
43	16.063	2202	2206	2211	rVB7	7589	12399	0.16%	0.012%
44	16.163	2218	2223	2230	rBV7	7043	14987	0.19%	0.015%
45	16.581	2290	2294	2298	rBV6	11015	18519	0.23%	0.018%
46	16.657	2303	2307	2311	rVV7	12181	16611	0.21%	0.017%
47	16.763	2320	2325	2330	rVB8	8861	14776	0.19%	0.015%
48	16.928	2351	2353	2358	rVB6	7627	8631	0.11%	0.009%
49	17.057	2368	2375	2380	rBV10	13557	30674	0.39%	0.031%
50	17.134	2383	2388	2392	rBV6	15701	24029	0.30%	0.024%
51	17.251	2401	2408	2414	rBV	2317426	3267320	41.37%	3.261%
52	17.345	2418	2424	2435	rVV	4468159	6445121	81.60%	6.432%
53	17.445	2439	2441	2444	rVB4	11747	10982	0.14%	0.011%
54	17.551	2454	2459	2463	rBV	237666	290951	3.68%	0.290%
55	17.604	2463	2468	2473	rVV	560485	637383	8.07%	0.636%
56	17.834	2504	2507	2512	rVB7	9662	13240	0.17%	0.013%
57	17.933	2518	2524	2527	rBV8	8600	16029	0.20%	0.016%
58	18.145	2555	2560	2571	rBV6	67715	142777	1.81%	0.142%
59	18.475	2610	2616	2623	rVB4	24756	41381	0.52%	0.041%
60	18.639	2641	2644	2649	rVB6	11178	13108	0.17%	0.013%
61	18.839	2673	2678	2681	rBV	344282	386208	4.89%	0.385%
62	18.881	2681	2685	2695	rVB	760997	874011	11.07%	0.872%
63	19.069	2715	2717	2721	rVB5	19084	21989	0.28%	0.022%
64	19.163	2729	2733	2737	rVB	59776	76131	0.96%	0.076%
65	19.233	2741	2745	2749	rBV2	51844	70119	0.89%	0.070%
66	19.386	2767	2771	2778	rBV4	78903	112937	1.43%	0.113%
67	19.445	2779	2781	2786	rVB6	20287	28672	0.36%	0.029%
68	19.586	2799	2805	2811	rBV	5458901	6928124	87.72%	6.914%
69	19.657	2811	2817	2823	rVB	327261	487328	6.17%	0.486%
70	19.839	2844	2848	2854	rVB	146626	173038	2.19%	0.173%
71	19.933	2859	2864	2867	rBV	3778057	3875565	49.07%	3.868%
72	19.969	2867	2870	2879	rVB	6793592	7898170	100.00%	7.882%
73	20.586	2971	2975	2979	rBV	233172	258559	3.27%	0.258%
74	20.898	3024	3028	3031	rBV	1282594	1421689	18.00%	1.419%
75	20.933	3031	3034	3047	rVB	2986959	3138818	39.74%	3.133%
76	21.351	3100	3105	3113	rBV	2635006	3218192	40.75%	3.212%

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77	21.422	3114	3117	3122	rVB2	137635	188511	2.39%	0.188%
78	21.810	3179	3183	3185	rBV	422599	522379	6.61%	0.521%
79	21.845	3185	3189	3201	rVB	1003827	1349148	17.08%	1.346%
80	22.022	3215	3219	3230	rVB	275682	443039	5.61%	0.442%
81	22.504	3298	3301	3306	rVB	148840	208215	2.64%	0.208%
82	22.698	3329	3334	3342	rVB	331696	659213	8.35%	0.658%
83	22.869	3359	3363	3366	rBV2	229550	376155	4.76%	0.375%
84	22.910	3366	3370	3383	rVB	576223	1043075	13.21%	1.041%
85	23.492	3464	3469	3480	rVB2	283540	724058	9.17%	0.723%
86	23.621	3484	3491	3501	rBV2	4002454	7814089	98.94%	7.799%
87	23.780	3511	3518	3532	rVB	1870397	3857944	48.85%	3.850%

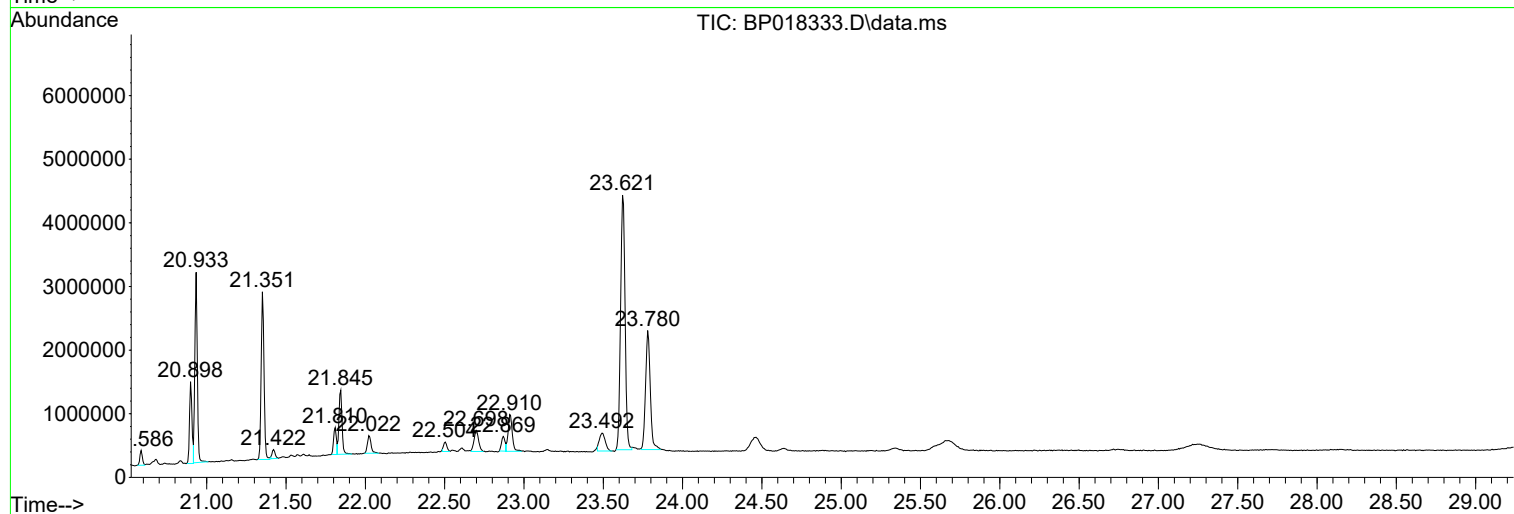
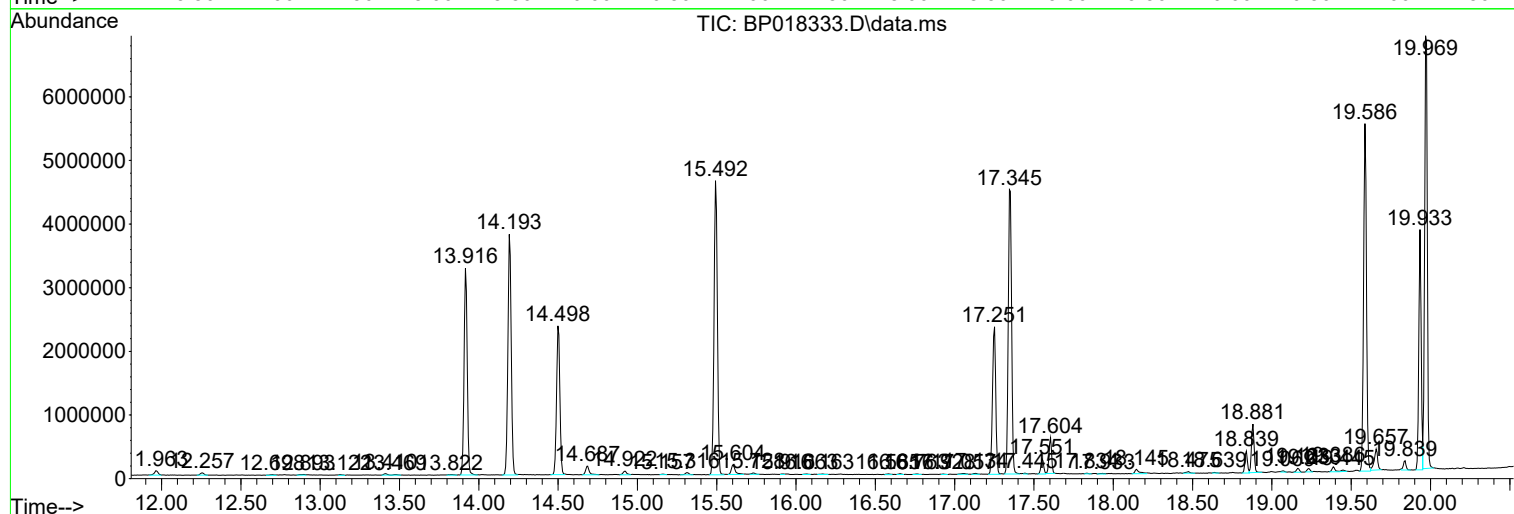
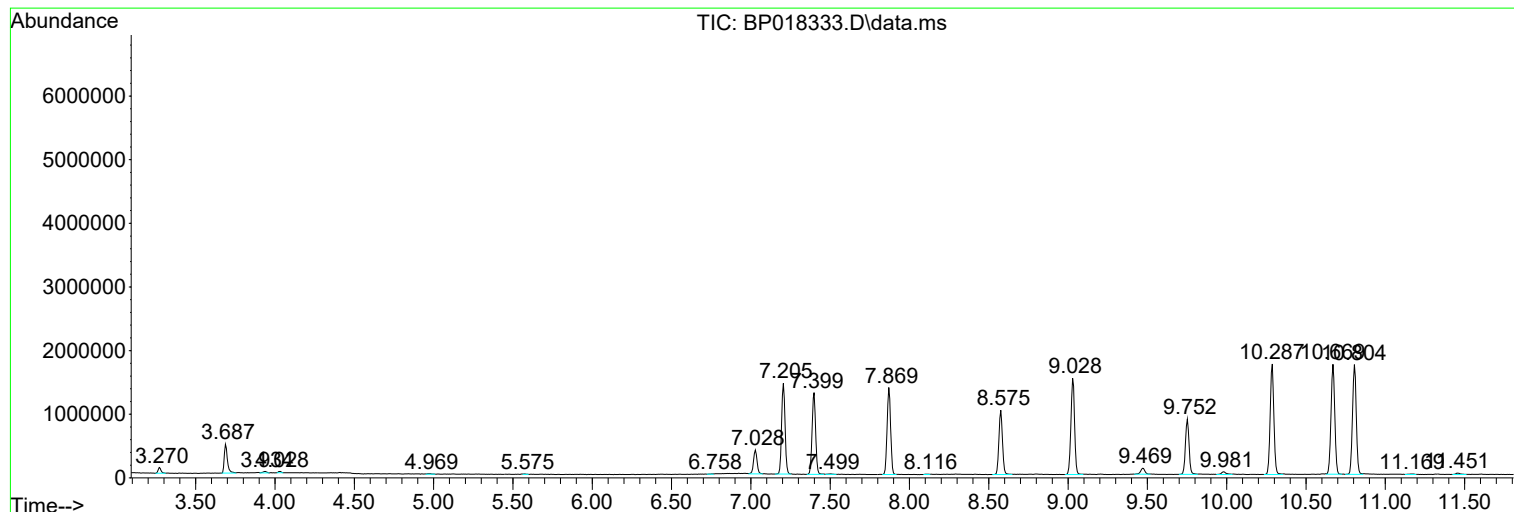
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112223.MA.M
Quant Title : SVOA CALIBRATION

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



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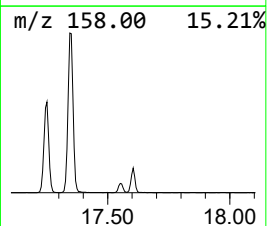
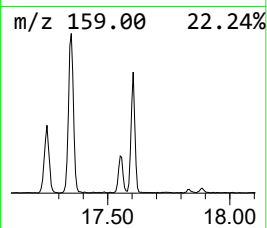
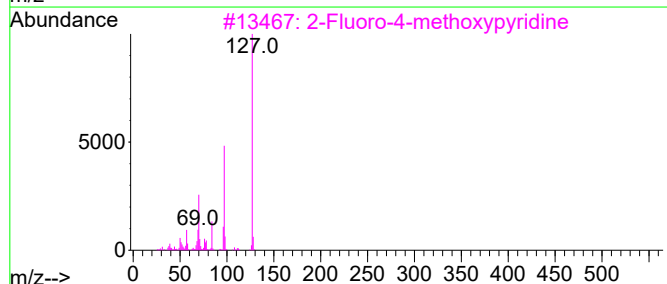
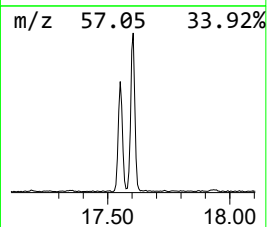
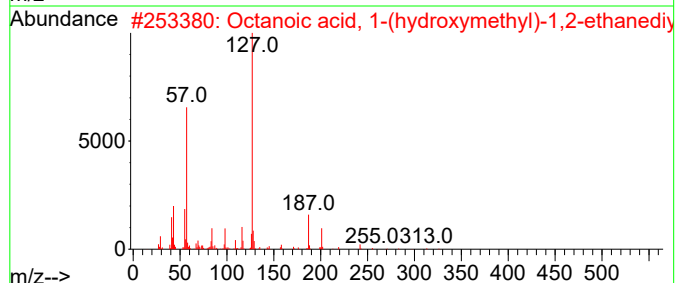
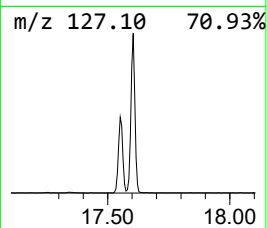
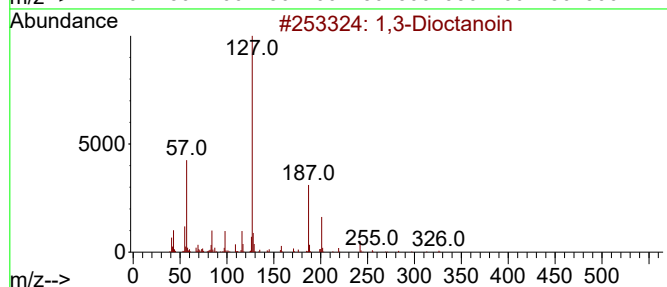
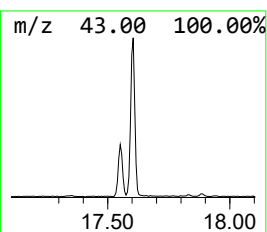
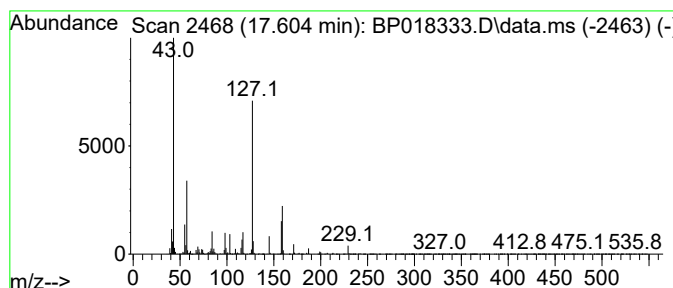
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112223.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 unknown-01 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.604	3.90 ng/ul	637383	Phenanthrene-d10	17.251

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1,3-Dioctanoin	344	C19H36O5	001429-66-9	38
2			Octanoic acid, 1-(hydroxymethyl)...	344	C19H36O5	060514-48-9	38
3			2-Fluoro-4-methoxypyridine	127	C6H6FNO	175965-83-0	35
4			2-Hydroxy-4-hydroxylaminopyrimidine	127	C4H5N3O2	020555-88-8	27
5			6-Azathymine	127	C4H5N3O2	000932-53-6	27



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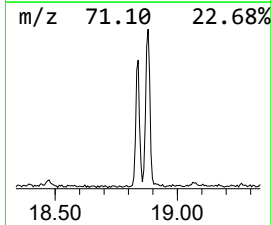
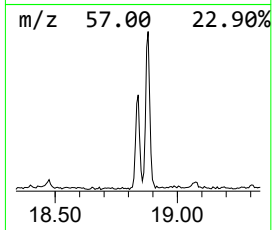
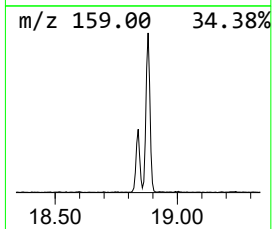
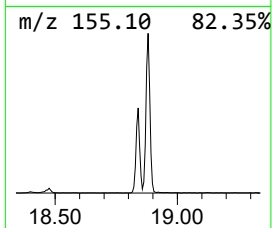
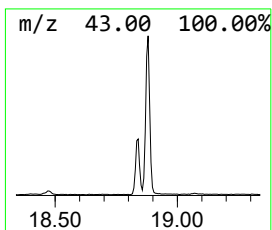
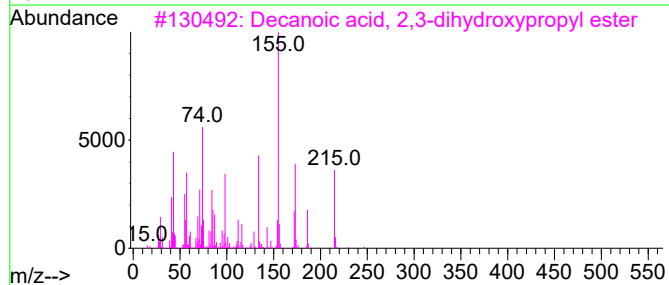
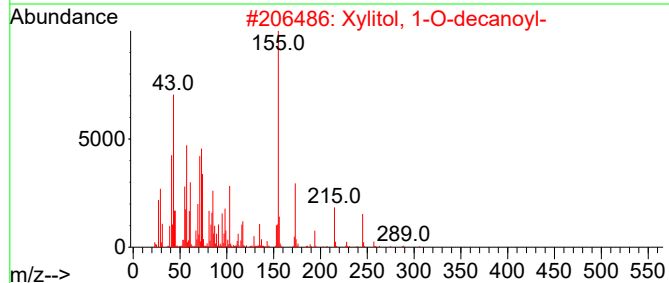
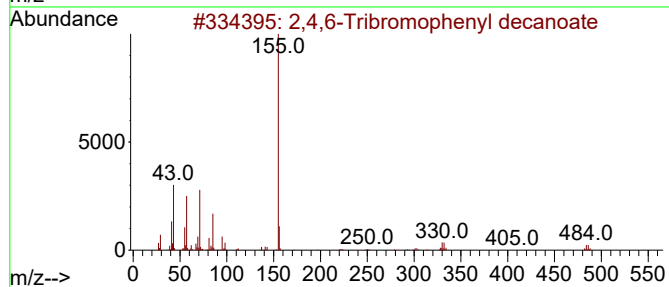
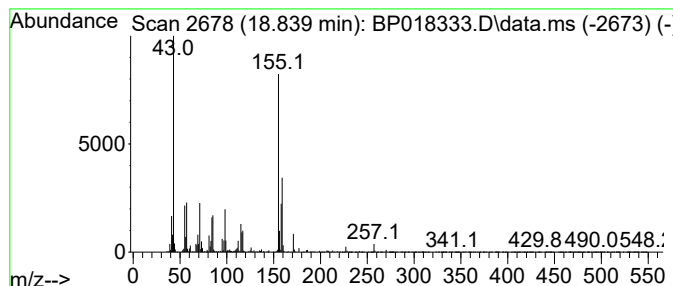
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112223.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 unknown-02 Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.839	2.36 ng/ul	386208	Phenanthrene-d10	17.251

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,4,6-Tribromophenyl decanoate	482	C16H21Br3O2	1000233-09-1	47		
2	Xylitol, 1-O-decanoyl-	306	C15H30O6	1000155-78-0	45		
3	Decanoic acid, 2,3-dihydroxyprop...	246	C13H26O4	002277-23-8	40		
4	Fumaric acid, butyl 3-heptyl ester	270	C15H26O4	1000348-68-0	38		
5	1,2-Cyclohexanedicarboxylic acid...	340	C20H36O4	1000339-45-2	38		



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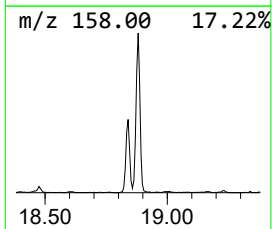
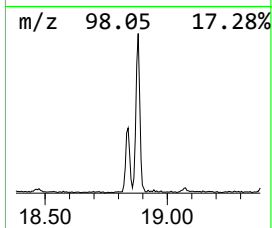
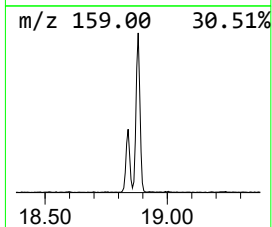
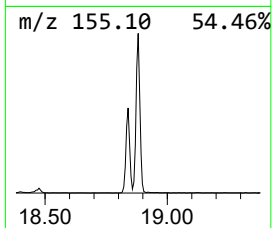
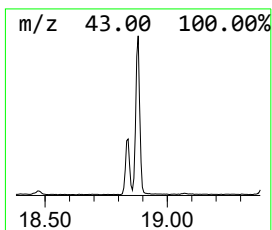
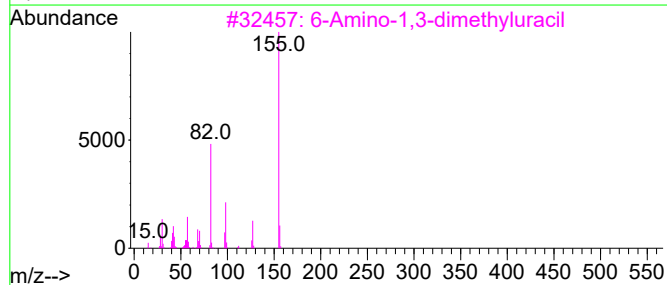
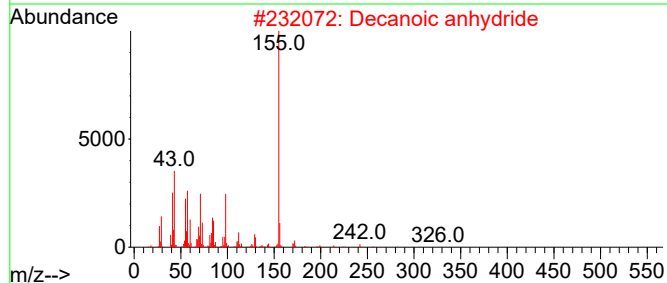
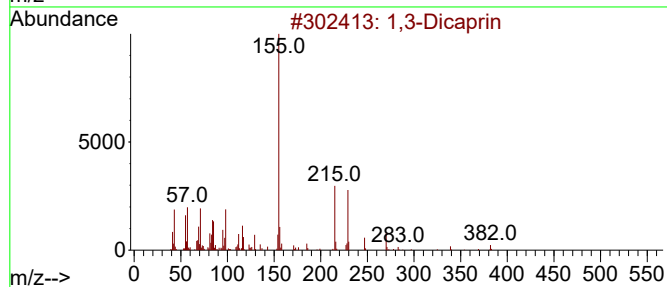
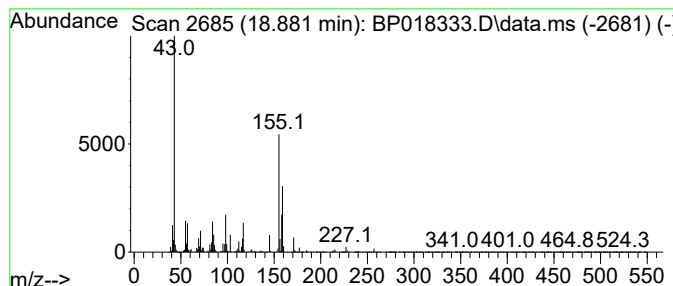
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112223.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 unknown-03 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.881	5.35 ng/ul	874011	Phenanthrene-d10	17.251

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1,3-Dicaprin	400	C23H44O5	017598-93-5	43
2			Decanoic anhydride	326	C20H38O3	002082-76-0	38
3			6-Amino-1,3-dimethyluracil	155	C6H9N3O2	006642-31-5	35
4			Decanoic anhydride	326	C20H38O3	002082-76-0	35
5			2,6-Difluoro-3-methylbenzoic aci...	354	C21H32F2O2	1000338-58-3	32



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DCKP7

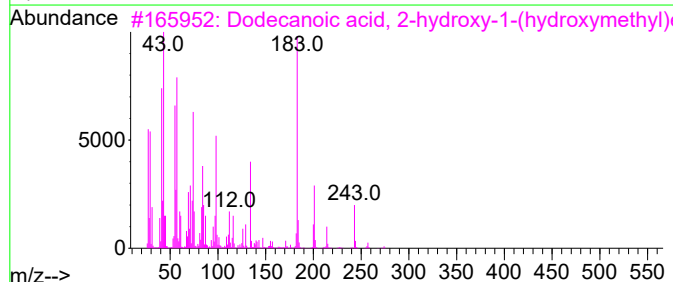
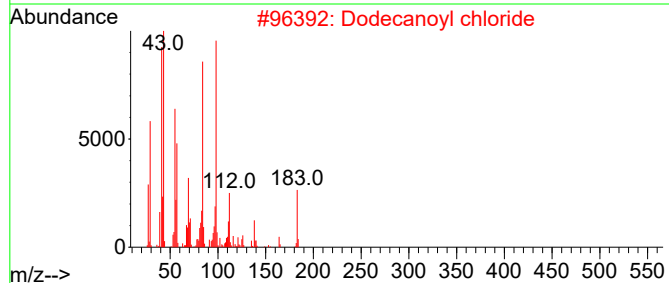
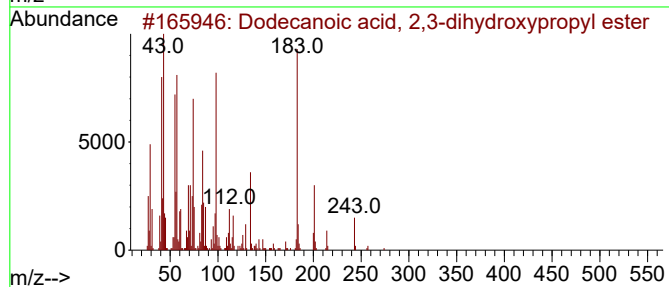
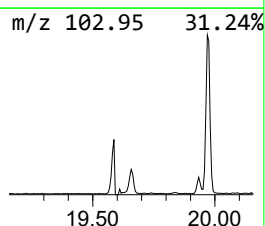
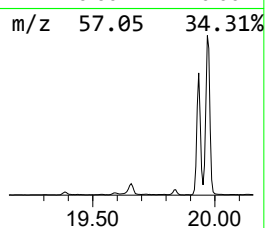
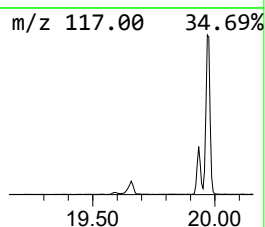
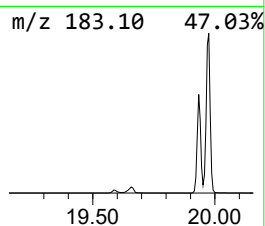
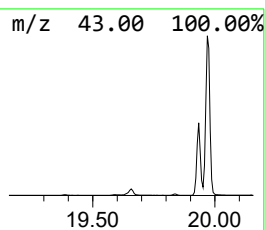
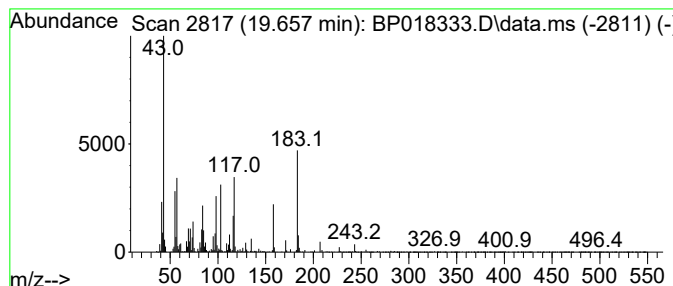
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 unknown-04 Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.657	3.03 ng/ul	487328	Chrysene-d12	21.351

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Dodecanoic acid, 2,3-dihydroxypr...	274	C15H30O4	000142-18-7	28
2			Dodecanoyl chloride	218	C12H23ClO	000112-16-3	16
3			Dodecanoic acid, 2-hydroxy-1-(hy...	274	C15H30O4	001678-45-1	10
4			9-Octadecenoic acid (Z)-, hexyl ...	366	C24H46O2	020290-84-0	9
5			Dodecanoic acid, ethenyl ester	226	C14H26O2	002146-71-6	9



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP112423\
Data File : BP018333.D
Acq On : 24 Nov 2023 18:26
Operator : MA/JU
Sample : 05264-09
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCKP7

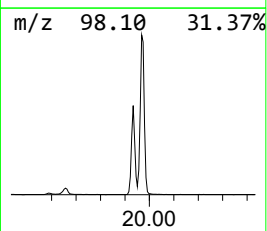
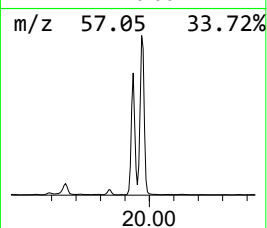
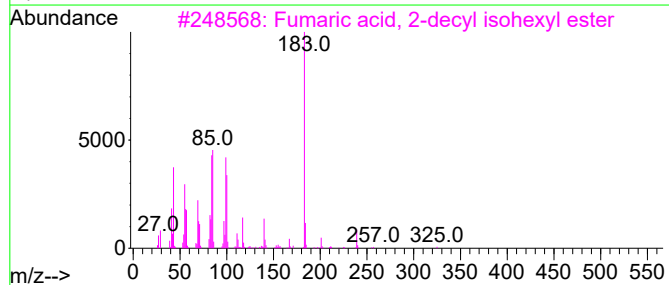
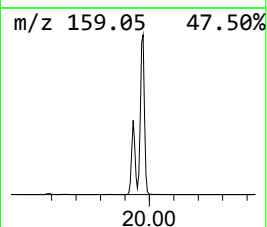
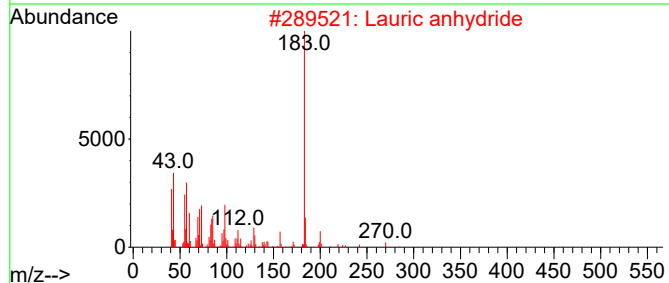
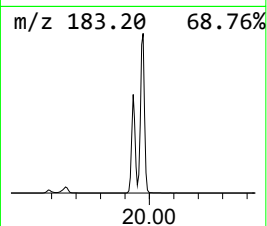
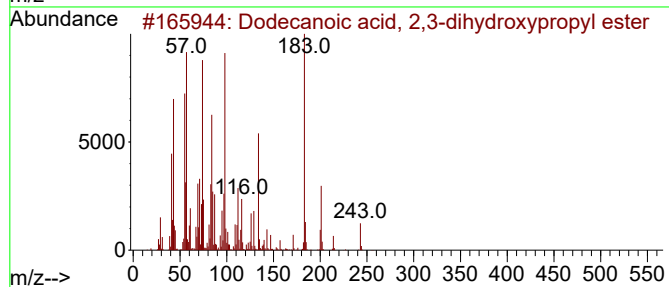
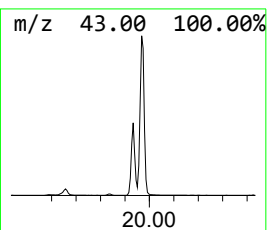
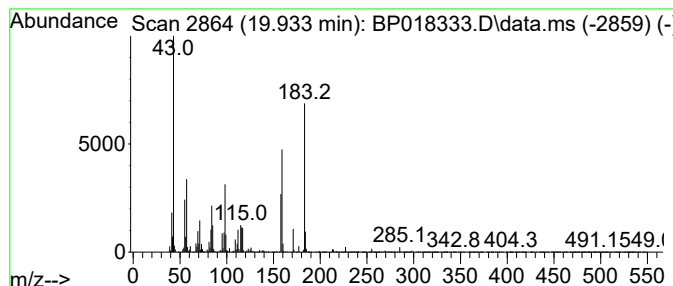
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 unknown-05 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.933	24.09 ng/ul	3875570	Chrysene-d12	21.351

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Dodecanoic acid, 2,3-dihydroxypr...	274	C15H30O4	000142-18-7	32
2			Lauric anhydride	382	C24H46O3	000645-66-9	27
3			Fumaric acid, 2-decyl isohexyl e...	340	C20H36O4	1000348-58-8	25
4			Fumaric acid, 2-decyl hexyl ester	340	C20H36O4	1000348-58-9	22
5			Fumaric acid, cyclobutyl isohexy...	254	C14H22O4	1000345-03-5	22



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP112423\
Data File : BP018333.D
Acq On : 24 Nov 2023 18:26
Operator : MA/JU
Sample : 05264-09
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCKP7

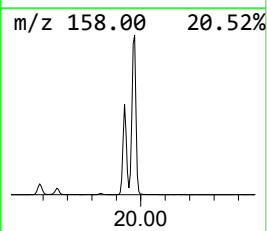
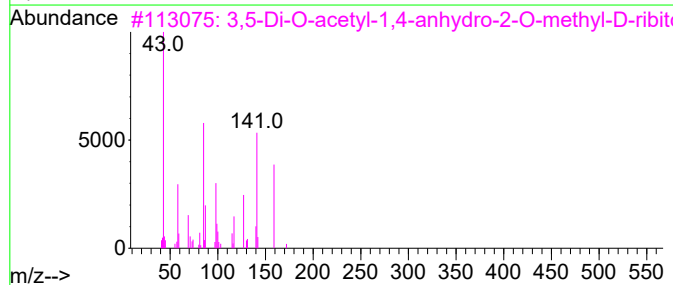
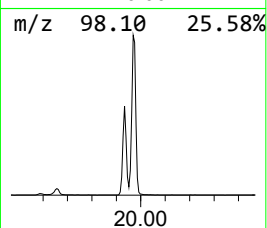
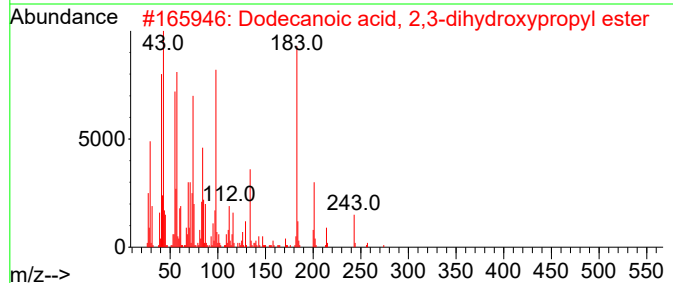
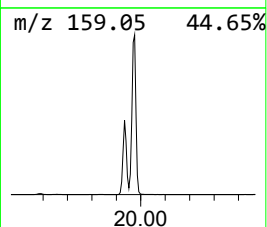
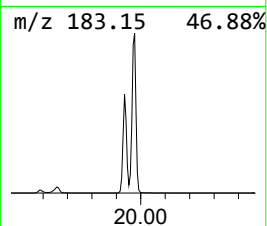
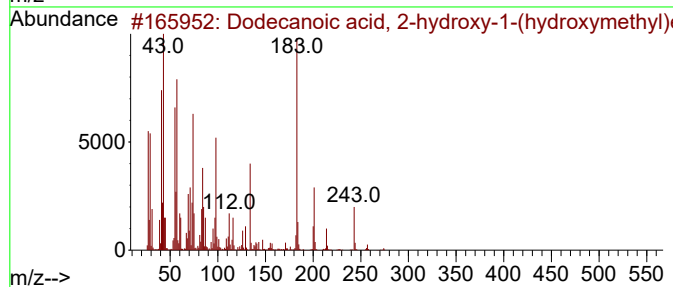
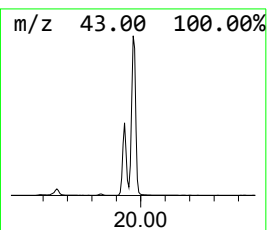
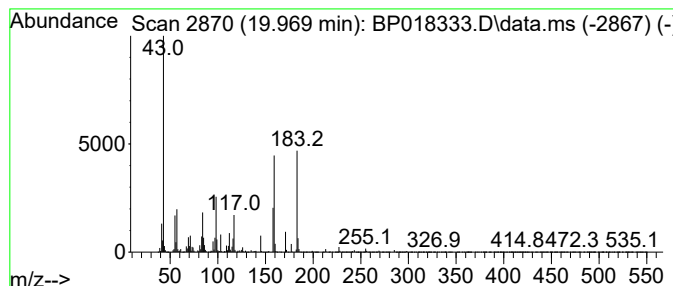
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 unknown-06 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.969	49.08 ng/ul	7898170	Chrysene-d12	21.351

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Dodecanoic acid, 2-hydroxy-1-(hy...	274	C15H30O4	001678-45-1	35
2			Dodecanoic acid, 2,3-dihydroxypr...	274	C15H30O4	000142-18-7	14
3			3,5-Di-O-acetyl-1,4-anhydro-2-O-...	232	C10H16O6	1010357-23-2	12
4			Eicosanoic acid, 2,3-bis(acetylo...	470	C27H50O6	055429-67-9	10
5			4,5,6-Trimethyltetrahydro-1,3-ox...	159	C7H13NO5	085333-98-8	9



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP112423\
Data File : BP018333.D
Acq On : 24 Nov 2023 18:26
Operator : MA/JU
Sample : 05264-09
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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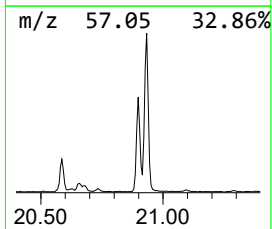
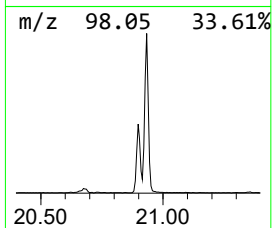
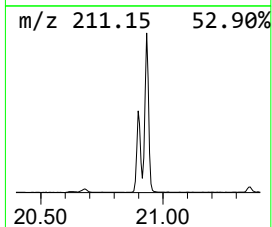
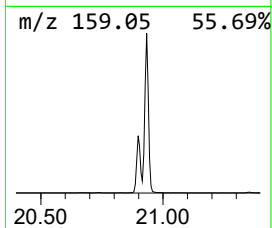
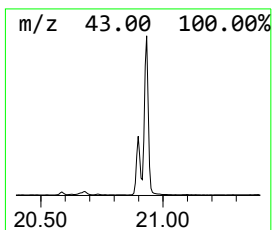
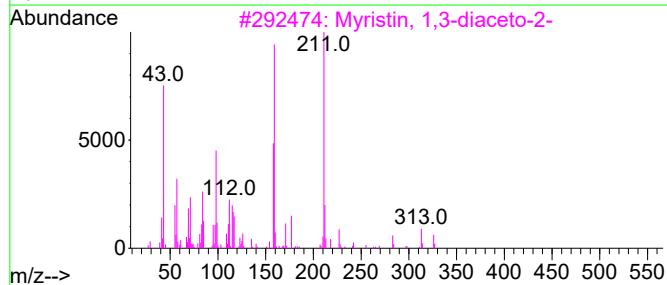
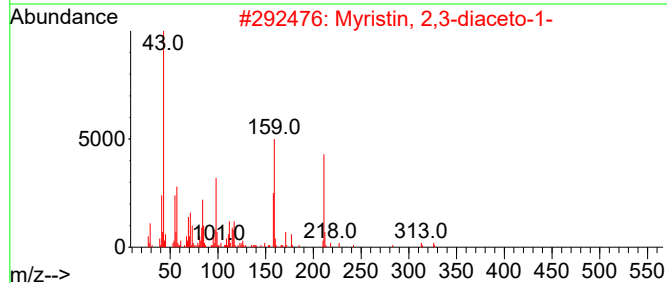
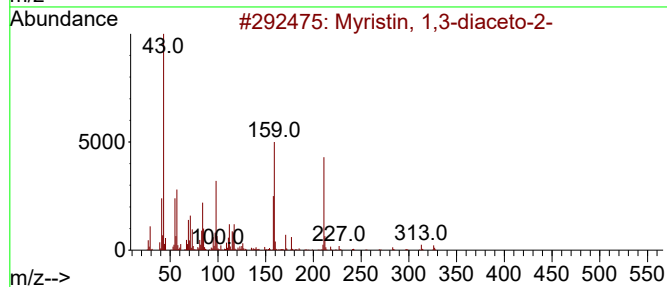
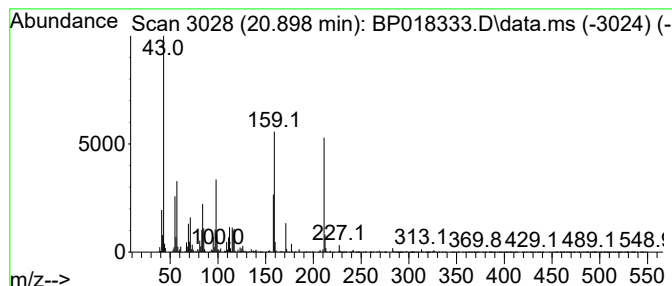
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 Myristin, 1,3-diaceto-2- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.898	8.84 ng/ul	1421690	Chrysene-d12	21.351

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Myristin, 1,3-diaceto-2-	386	C21H38O6	014290-23-4	91		
2	Myristin, 2,3-diaceto-1-	386	C21H38O6	014473-55-3	91		
3	Myristin, 1,3-diaceto-2-	386	C21H38O6	014290-23-4	72		
4	Myristin, 2,3-diaceto-1-	386	C21H38O6	014473-55-3	64		
5	Hexadecanoic acid, 2,3-bis(acety...	414	C23H42O6	055268-70-7	52		



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP112423\
Data File : BP018333.D
Acq On : 24 Nov 2023 18:26
Operator : MA/JU
Sample : 05264-09
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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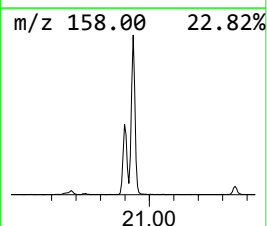
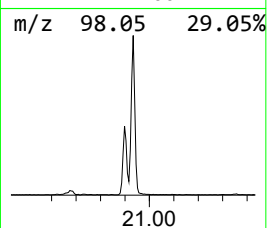
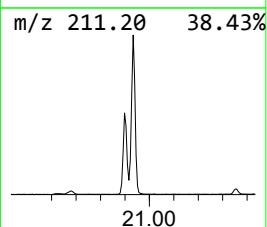
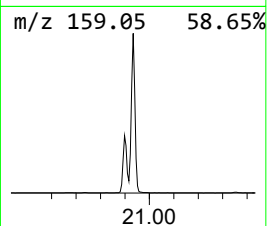
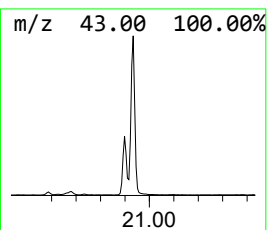
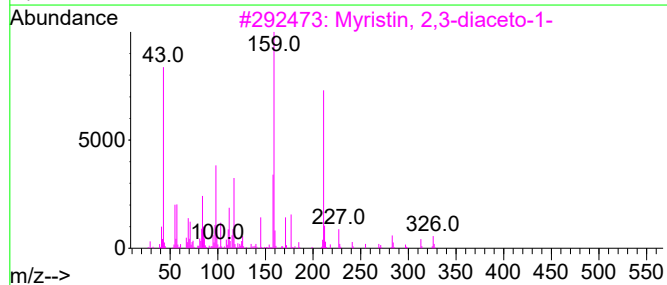
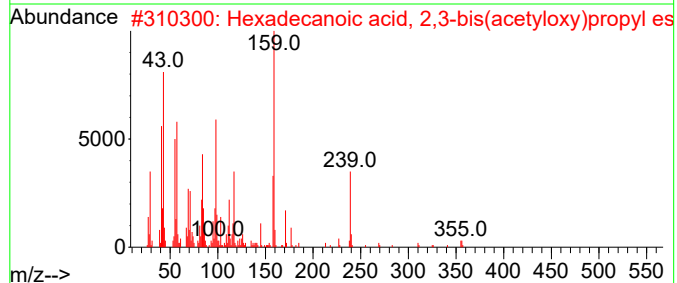
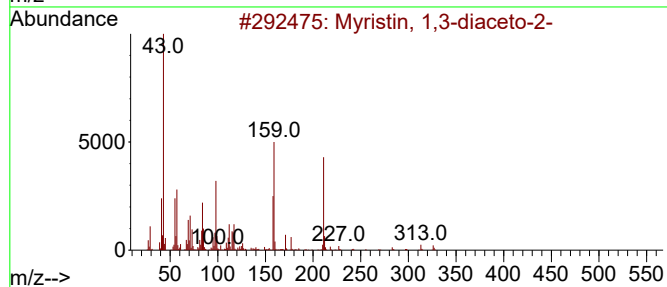
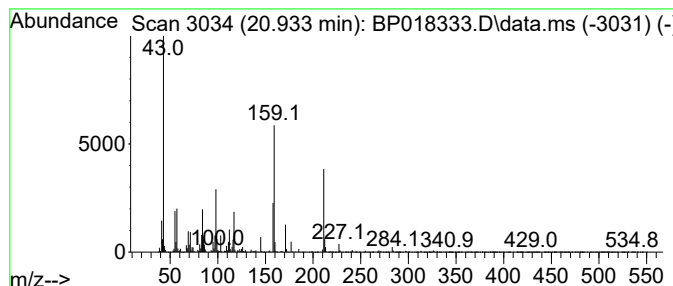
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 unknown-07 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.933	19.51 ng/ul	3138820	Chrysene-d12	21.351

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Myristin, 1,3-diaceto-2-	386	C21H38O6	014290-23-4	74		
2	Hexadecanoic acid, 2,3-bis(acety...	414	C23H42O6	055268-70-7	49		
3	Myristin, 2,3-diaceto-1-	386	C21H38O6	014473-55-3	45		
4	Myristin, 2,3-diaceto-1-	386	C21H38O6	014473-55-3	38		
5	2,3-Dihydroxypropyl 12-methyltri...	386	C21H38O6	1000488-66-8	38		



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP112423\
Data File : BP018333.D
Acq On : 24 Nov 2023 18:26
Operator : MA/JU
Sample : 05264-09
Misc :
ALS Vial : 14 Sample Multiplier: 1

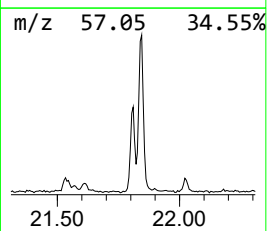
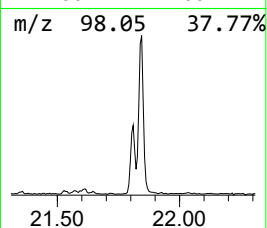
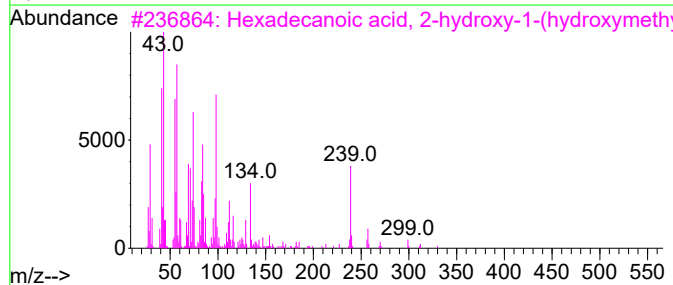
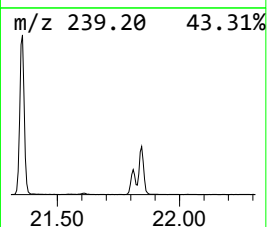
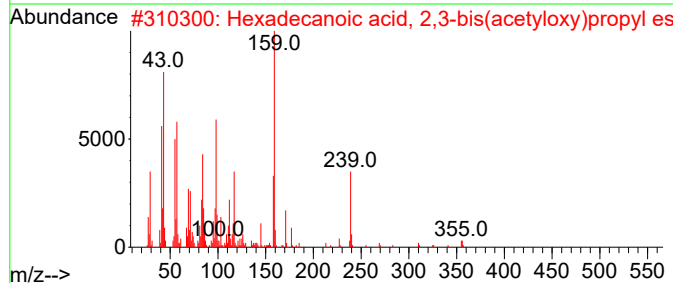
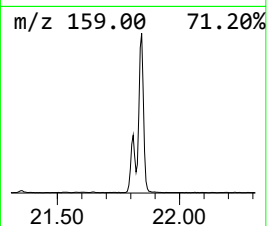
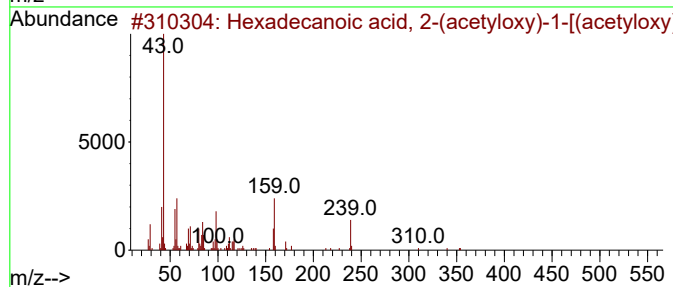
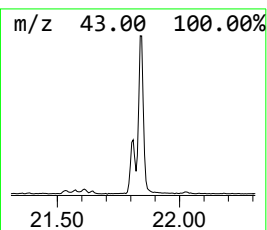
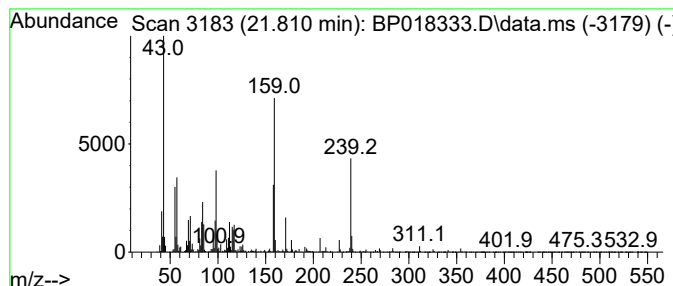
Instrument :
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112223.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 9 Hexadecanoic acid, 2-(acety... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.	
21.810	3.25 ng/ul	522379	Chrysene-d12	21.351	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexadecanoic acid, 2-(acetyloxy)...	414	C23H42O6	055268-69-4	91
2	Hexadecanoic acid, 2,3-bis(acety...	414	C23H42O6	055268-70-7	90
3	Hexadecanoic acid, 2-hydroxy-1-(...	330	C19H38O4	023470-00-0	30
4	Eicosanoic acid, 2-(acetyloxy)-1...	470	C27H50O6	055429-68-0	27
5	Tetradecyl nonanoate	354	C23H46O2	072934-14-6	14



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP112423\
Data File : BP018333.D
Acq On : 24 Nov 2023 18:26
Operator : MA/JU
Sample : 05264-09
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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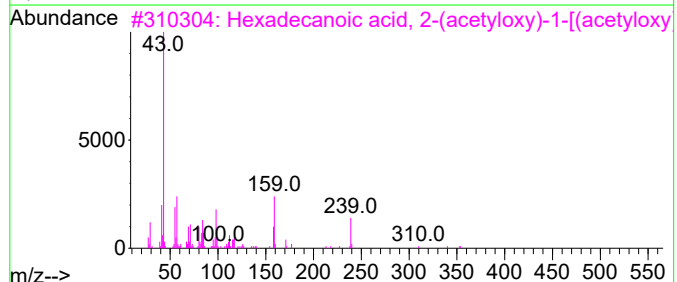
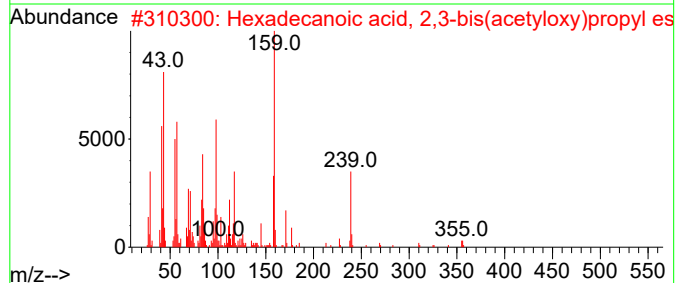
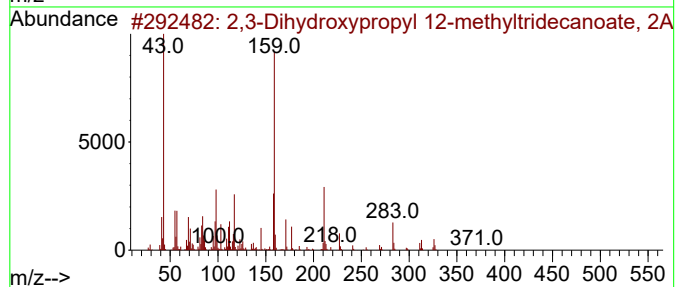
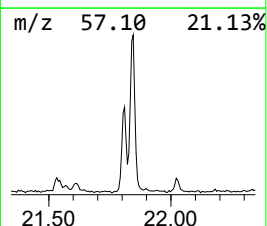
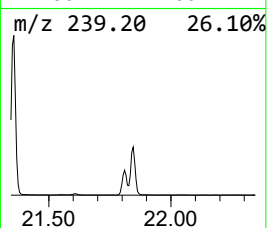
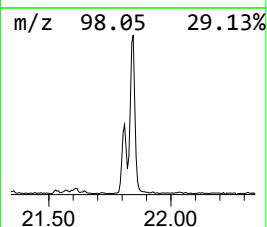
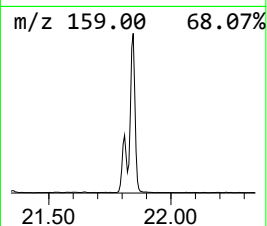
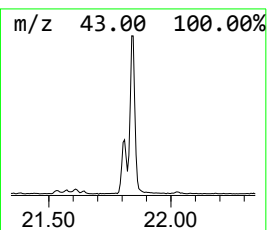
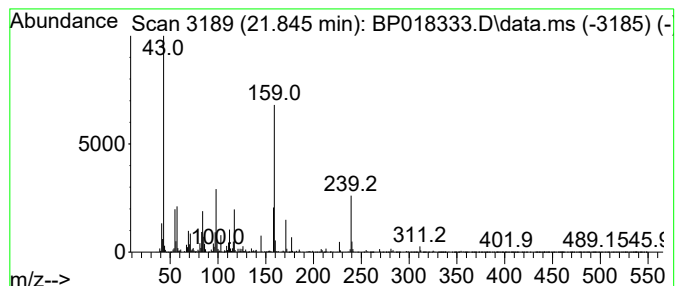
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 2,3-Dihydroxypropyl 12-meth... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.845	8.38 ng/ul	1349150	Chrysene-d12	21.351

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,3-Dihydroxypropyl 12-methyltri...	386	C21H38O6	1000488-66-8	53		
2	Hexadecanoic acid, 2,3-bis(acety...	414	C23H42O6	055268-70-7	52		
3	Hexadecanoic acid, 2-(acetyloxy)...	414	C23H42O6	055268-69-4	43		
4	3-((12-Acetoxyoctadecanoyl)oxy)p...	500	C27H48O8	330198-91-9	16		
5	2,3,4-Trifluorobenzoic acid, 3-t...	358	C20H29F3O2	1000280-61-8	16		



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP112423\
Data File : BP018333.D
Acq On : 24 Nov 2023 18:26
Operator : MA/JU
Sample : 05264-09
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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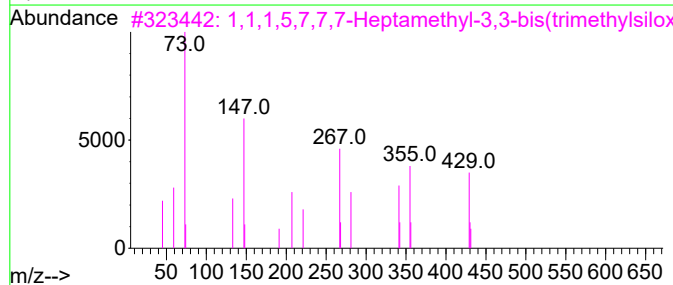
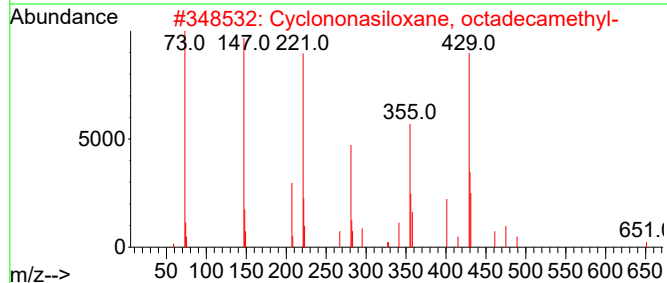
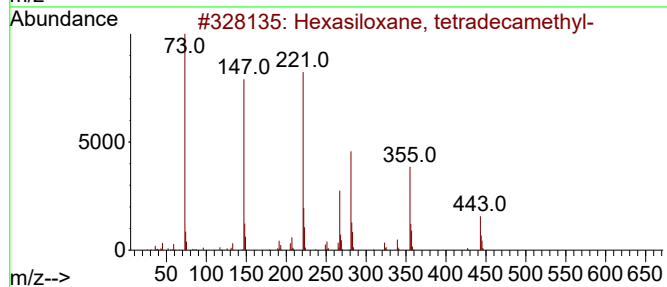
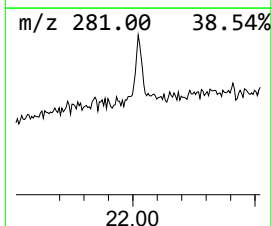
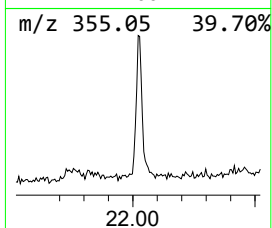
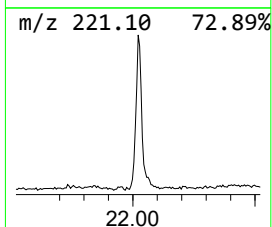
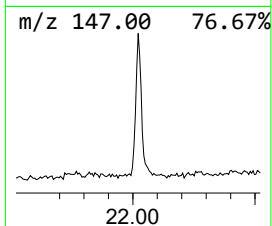
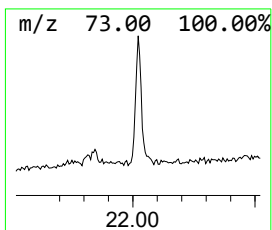
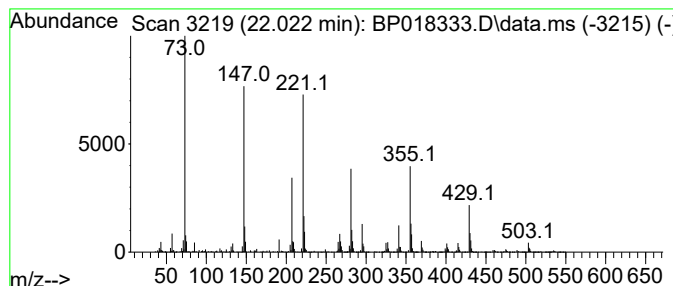
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Quant Title : SVOA CALIBRATION

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

Peak Number 11 Hexasiloxane, tetradecamethyl- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.022	2.75 ng/ul	443039	Chrysene-d12	21.351

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexasiloxane, tetradecamethyl-	458	C14H42O5Si6	000107-52-8	74
2			Cyclononasiloxane, octadecamethyl-	666	C18H54O9Si9	000556-71-8	74
3			1,1,1,5,7,7,7-Heptamethyl-3,3-bi...	444	C13H40O5Si6	038147-00-1	59
4			Cyclodecasiloxane, eicosamethyl-	740	C20H60O10Si10	018772-36-6	45
5			N-Benzyl-N-ethyl-p-isopropylbenz...	281	C19H23NO	015089-22-2	43



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP112423\
Data File : BP018333.D
Acq On : 24 Nov 2023 18:26
Operator : MA/JU
Sample : 05264-09
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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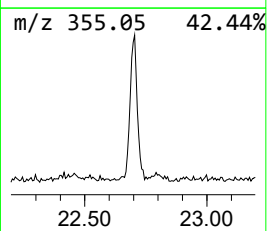
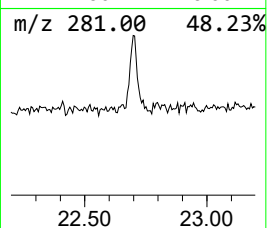
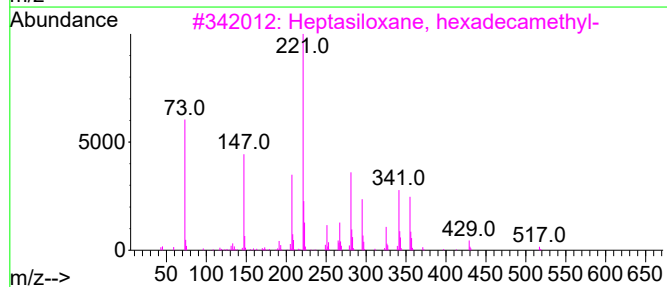
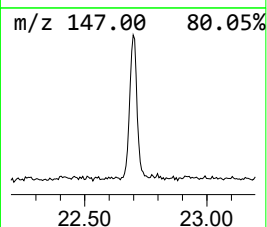
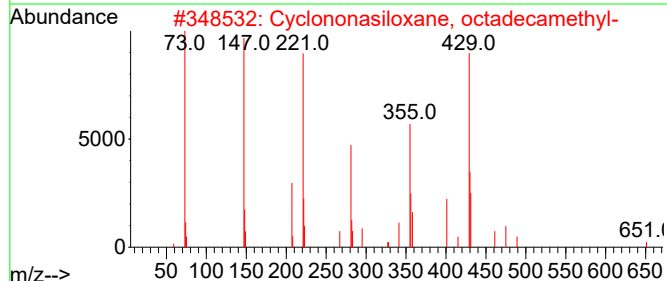
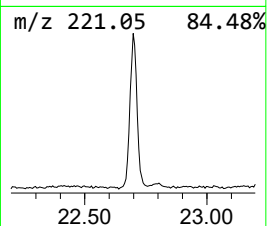
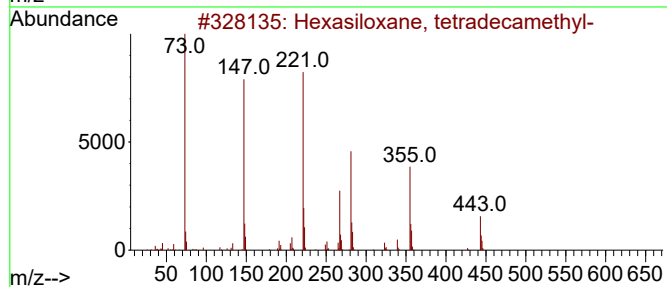
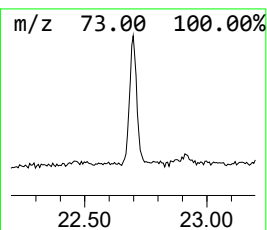
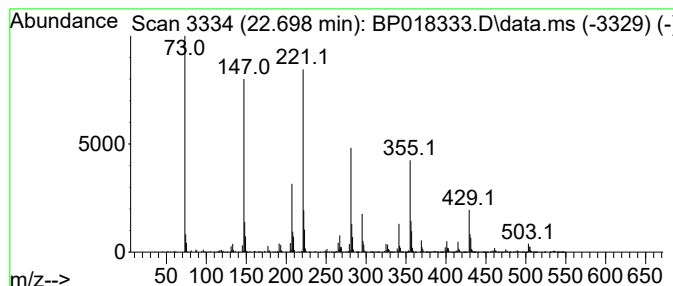
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Quant Title : SVOA CALIBRATION

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

Peak Number 12 Cyclononasiloxane, octadeca... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.698	3.42 ng/ul	659213	Perylene-d12	23.780

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexasiloxane, tetradecamethyl-	458	C14H42O5Si6	000107-52-8	87
2			Cyclononasiloxane, octadecamethyl-	666	C18H54O9Si9	000556-71-8	78
3			Heptasiloxane, hexadecamethyl-	532	C16H48O6Si7	000541-01-5	45
4			Pentasiloxane, dodecamethyl-	384	C12H36O4Si5	000141-63-9	30
5			N-Benzyl-N-ethyl-p-isopropylbenz...	281	C19H23NO	015089-22-2	27



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP112423\
Data File : BP018333.D
Acq On : 24 Nov 2023 18:26
Operator : MA/JU
Sample : 05264-09
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCKP7

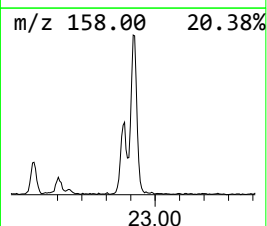
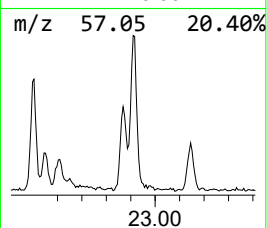
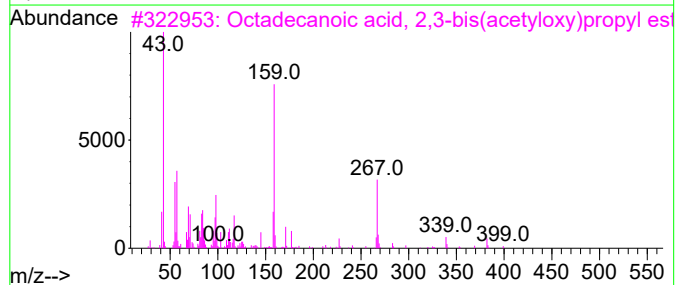
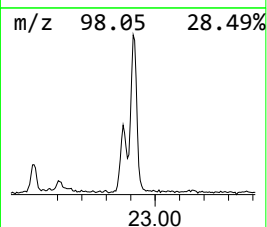
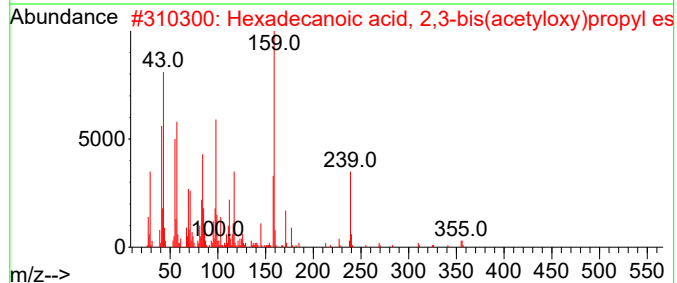
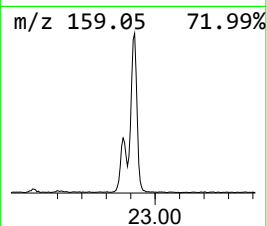
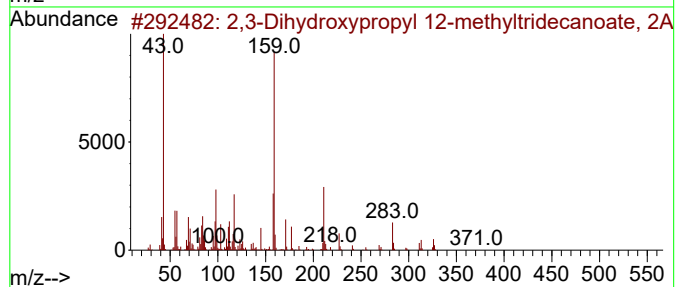
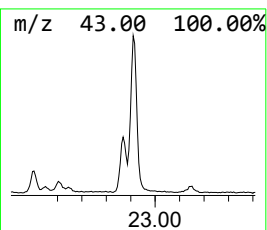
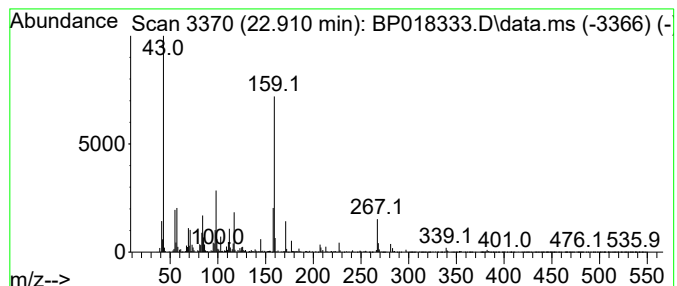
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Quant Title : SVOA CALIBRATION

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

Peak Number 13 unknown-08 Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.910	5.41 ng/ul	1043080	Perylene-d12	23.780

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,3-Dihydroxypropyl 12-methyltri...	386	C21H38O6	1000488-66-8	72		
2	Hexadecanoic acid, 2,3-bis(acety...	414	C23H42O6	055268-70-7	49		
3	Octadecanoic acid, 2,3-bis(acety...	442	C25H46O6	033599-07-4	45		
4	Octadecanoic acid, 2-(acetyloxy)...	442	C25H46O6	055401-62-2	37		
5	2,3,4-Trifluorobenzoic acid, 3-t...	372	C21H31F3O2	1000280-61-9	27		



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP112423\
Data File : BP018333.D
Acq On : 24 Nov 2023 18:26
Operator : MA/JU
Sample : 05264-09
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCKP7

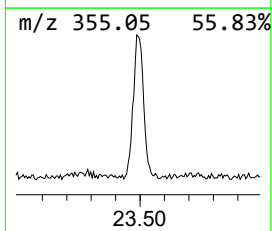
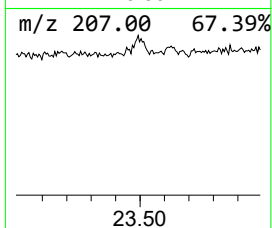
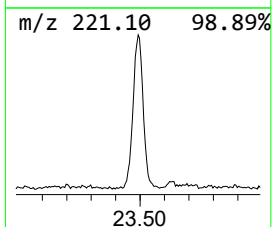
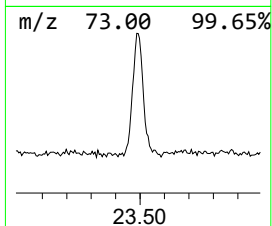
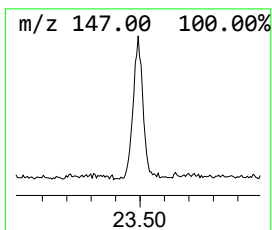
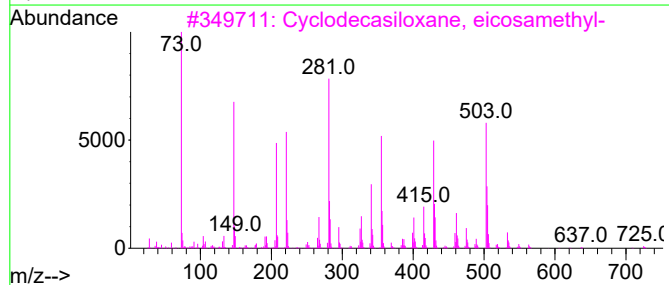
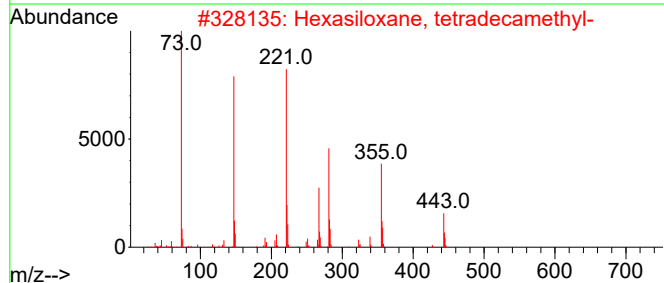
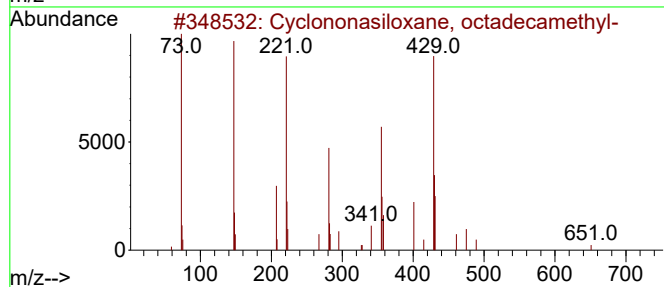
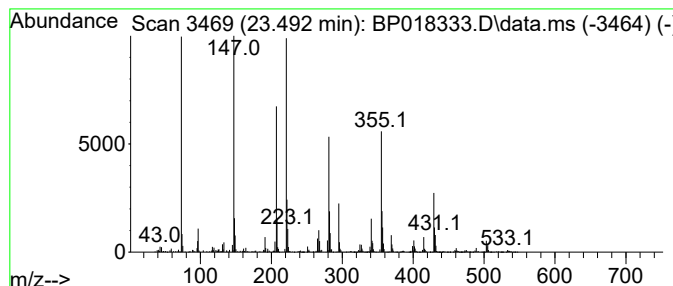
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Quant Title : SVOA CALIBRATION

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

Peak Number 14 unknown-09 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.492	3.75 ng/ul	724058	Perylene-d12	23.780

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclononasiloxane, octadecamethyl-	666	C18H54O9Si9	000556-71-8	86
2			Hexasiloxane, tetradecamethyl-	458	C14H42O5Si6	000107-52-8	72
3			Cyclodecasiloxane, eicosamethyl-	740	C20H60O10Si10	018772-36-6	42
4			Mercaptoacetic acid, 2TMS deriva...	236	C8H20O2SSi2	006398-62-5	32
5			Heptasiloxane, hexadecamethyl-	532	C16H48O6Si7	000541-01-5	32



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP112423\
 Data File : BP018333.D
 Acq On : 24 Nov 2023 18:26
 Operator : MA/JU
 Sample : 05264-09
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 DCKP7

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112223.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
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unknown-02	18.839	2.4	ng/ul	386208	4	17.251	3267320	20.0
unknown-03	18.881	5.3	ng/ul	874011	4	17.251	3267320	20.0
unknown-04	19.657	3.0	ng/ul	487328	5	21.351	3218190	20.0
unknown-05	19.933	24.1	ng/ul	3875570	5	21.351	3218190	20.0
unknown-06	19.969	49.1	ng/ul	7898170	5	21.351	3218190	20.0
Myristin, 1,3-d...	20.898	8.8	ng/ul	1421690	5	21.351	3218190	20.0
unknown-07	20.933	19.5	ng/ul	3138820	5	21.351	3218190	20.0
Hexadecanoic ac...	21.810	3.3	ng/ul	522379	5	21.351	3218190	20.0
2,3-Dihydroxypr...	21.845	8.4	ng/ul	1349150	5	21.351	3218190	20.0
Hexasiloxane, t...	22.022	2.8	ng/ul	443039	5	21.351	3218190	20.0
Cyclononasiloxa...	22.698	3.4	ng/ul	659213	6	23.780	3857940	20.0
unknown-08	22.910	5.4	ng/ul	1043080	6	23.780	3857940	20.0
unknown-09	23.492	3.8	ng/ul	724058	6	23.780	3857940	20.0