

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121922\
 Data File : BP013146.D
 Acq On : 20 Dec 2022 02:18
 Operator : CG/JU
 Sample : N6006-06
 Misc :
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 DBXV9

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-BP120722.M
 Title : SVOA CALIBRATION

Signal : TIC: BP013146.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.334	22	25	36	rVB	199088	284231	1.61%	0.113%
2	3.770	95	99	112	rBV	1027318	1574938	8.92%	0.626%
3	3.876	113	117	122	rVV	127359	173585	0.98%	0.069%
4	3.999	134	138	143	rVB	81082	114984	0.65%	0.046%
5	4.099	150	155	159	rBV	44279	64248	0.36%	0.026%
6	4.664	247	251	257	rVB	36841	55235	0.31%	0.022%
7	5.040	308	315	325	rVB2	255989	416077	2.36%	0.165%
8	5.275	350	355	362	rBV4	26380	48225	0.27%	0.019%
9	6.628	579	585	590	rVB	62163	93744	0.53%	0.037%
10	7.111	658	667	682	rBV	3045968	5010616	28.39%	1.993%
11	7.281	689	696	709	rBV	2857755	4361065	24.71%	1.734%
12	7.481	722	730	740	rBV	3261719	5230275	29.63%	2.080%
13	7.958	803	811	818	rBV	2114717	3351231	18.98%	1.333%
14	8.087	828	833	842	rVB	88019	137664	0.78%	0.055%
15	8.663	923	931	941	rBV	3092859	4963823	28.12%	1.974%
16	8.787	946	952	963	rVB	169630	293892	1.66%	0.117%
17	9.122	1002	1009	1017	rBV	3341648	5440964	30.82%	2.164%
18	9.528	1068	1078	1084	rBV2	83298	140773	0.80%	0.056%
19	9.846	1124	1132	1148	rBV	2770962	4776752	27.06%	1.900%
20	9.987	1152	1156	1163	rVV2	30950	68756	0.39%	0.027%
21	10.063	1166	1169	1178	rVB7	31057	66649	0.38%	0.027%
22	10.387	1216	1224	1234	rBV	4284004	7005299	39.69%	2.786%
23	10.769	1278	1289	1296	rBV	3030841	5040741	28.56%	2.004%
24	10.905	1304	1312	1338	rVB	2057523	3834363	21.72%	1.525%
25	11.299	1374	1379	1390	rVB8	29944	63755	0.36%	0.025%
26	11.552	1415	1422	1429	rBV4	35795	75099	0.43%	0.030%
27	12.175	1523	1528	1537	rVB	49379	86995	0.49%	0.035%
28	12.287	1544	1547	1553	rVV	44107	70823	0.40%	0.028%
29	12.352	1553	1558	1564	rVB2	76540	121134	0.69%	0.048%
30	13.334	1719	1725	1728	rBV7	28617	54455	0.31%	0.022%
31	13.410	1734	1738	1745	rVB	95597	127531	0.72%	0.051%
32	13.481	1745	1750	1754	rBV5	40743	72279	0.41%	0.029%
33	13.916	1819	1824	1830	rBV3	121260	181990	1.03%	0.072%
34	14.004	1830	1839	1845	rBV	7795178	10644805	60.30%	4.233%
35	14.122	1854	1859	1863	rVB6	40868	62113	0.35%	0.025%
36	14.293	1881	1888	1898	rVV	8606153	12490145	70.76%	4.967%

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Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
 Title : SVOA CALIBRATION

37	14.481	1916	1920	1926	rVB2	47549	66676	0.38%	0.027%
38	14.598	1929	1940	1946	rBV	4759146	6951191	39.38%	2.764%
39	14.657	1946	1950	1958	rVB	99497	162207	0.92%	0.065%
40	14.793	1965	1973	1990	rBV	2878033	4556825	25.81%	1.812%
41	14.946	1995	1999	2004	rVV2	118776	161674	0.92%	0.064%
42	15.004	2004	2009	2015	rVB5	101527	187776	1.06%	0.075%
43	15.134	2026	2031	2035	rBV3	78532	103321	0.59%	0.041%
44	15.375	2068	2072	2075	rBV5	27073	46924	0.27%	0.019%
45	15.593	2101	2109	2115	rBV	11482782	16337102	92.55%	6.497%
46	15.645	2115	2118	2123	rVV	174189	228018	1.29%	0.091%
47	15.710	2123	2129	2136	rVV	3644837	5012940	28.40%	1.993%
48	15.840	2144	2151	2156	rVB3	133860	232570	1.32%	0.092%
49	15.957	2165	2171	2181	rVB	1107792	1501859	8.51%	0.597%
50	16.163	2202	2206	2212	rBV8	25152	46554	0.26%	0.019%
51	16.298	2225	2229	2238	rVB10	43996	106787	0.60%	0.042%
52	16.481	2256	2260	2265	rVV2	36538	55976	0.32%	0.022%
53	16.657	2278	2290	2295	rBV5	42174	121760	0.69%	0.048%
54	16.887	2325	2329	2332	rBV2	53668	67345	0.38%	0.027%
55	17.004	2345	2349	2355	rBV2	45039	77449	0.44%	0.031%
56	17.098	2359	2365	2368	rVB5	32050	58442	0.33%	0.023%
57	17.169	2368	2377	2383	rVB2	77298	161812	0.92%	0.064%
58	17.234	2384	2388	2396	rBV4	46078	75135	0.43%	0.030%
59	17.298	2396	2399	2402	rBV4	36223	45326	0.26%	0.018%
60	17.351	2402	2408	2412	rVV	5684838	8213385	46.53%	3.266%
61	17.392	2412	2415	2420	rVV	1523759	2163605	12.26%	0.860%
62	17.451	2420	2425	2435	rVV	10441929	15892949	90.03%	6.320%
63	17.534	2435	2439	2448	rVB3	250628	446113	2.53%	0.177%
64	17.728	2469	2472	2474	rVV3	46005	64244	0.36%	0.026%
65	17.757	2474	2477	2481	rVB2	52345	69514	0.39%	0.028%
66	18.110	2533	2537	2542	rVB3	54552	73449	0.42%	0.029%
67	18.169	2543	2547	2551	rBV	175026	241359	1.37%	0.096%
68	18.228	2551	2557	2562	rBV	3761274	4979135	28.21%	1.980%
69	18.410	2582	2588	2596	rBV4	437433	827148	4.69%	0.329%
70	18.469	2596	2598	2602	rVB	78309	81564	0.46%	0.032%
71	18.504	2602	2604	2608	rBV3	39671	50385	0.29%	0.020%
72	18.551	2610	2612	2617	rVB4	58062	70560	0.40%	0.028%
73	18.634	2622	2626	2633	rVB	309738	365506	2.07%	0.145%
74	18.698	2633	2637	2640	rBV2	203142	277978	1.57%	0.111%
75	18.734	2640	2643	2648	rVV4	131037	220141	1.25%	0.088%
76	18.786	2648	2652	2660	rVB	318536	415797	2.36%	0.165%

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77	18.934	2674	2677	2681	rVB3	58775	70519	0.40%	0.028%
78	19.104	2701	2706	2713	rBV4	169905	480628	2.72%	0.191%
79	19.204	2720	2723	2726	rBV2	81251	115788	0.66%	0.046%
80	19.257	2729	2732	2738	rVB2	220136	318261	1.80%	0.127%
81	19.357	2738	2749	2761	rBV	3793022	7649144	43.33%	3.042%
82	19.457	2761	2766	2779	rVV	1187166	1780260	10.09%	0.708%
83	19.592	2787	2789	2797	rVB2	120260	179870	1.02%	0.072%
84	19.686	2799	2805	2808	rBV	12116898	17264076	97.80%	6.865%
85	19.833	2825	2830	2835	rBV	2075634	2326476	13.18%	0.925%
86	19.886	2836	2839	2844	rVB	117098	156952	0.89%	0.062%
87	20.186	2884	2890	2896	rVB2	543416	1044248	5.92%	0.415%
88	20.286	2905	2907	2911	rVB2	168609	179945	1.02%	0.072%
89	21.128	3046	3050	3055	rBV2	2831079	3758133	21.29%	1.494%
90	21.439	3096	3103	3107	rBV2	4776562	7882265	44.65%	3.134%
91	21.475	3107	3109	3115	rVB	763471	974893	5.52%	0.388%
92	22.045	3203	3206	3208	rBV2	465840	619106	3.51%	0.246%
93	22.069	3208	3210	3220	rVB2	867143	1279782	7.25%	0.509%
94	23.133	3386	3391	3411	rVB2	2537746	6715649	38.04%	2.671%
95	23.757	3491	3497	3512	rVB2	6000144	13664813	77.41%	5.434%
96	23.916	3518	3524	3532	rBV2	2842473	5844264	33.11%	2.324%
97	24.539	3624	3630	3639	rVB	2184898	4730340	26.80%	1.881%
98	26.457	3948	3956	3976	rVB2	1547946	5751421	32.58%	2.287%
99	27.463	4122	4127	4141	rVB3	1149746	3616991	20.49%	1.438%
100	28.515	4297	4306	4330	rVB4	4150633	17652012	100.00%	7.019%

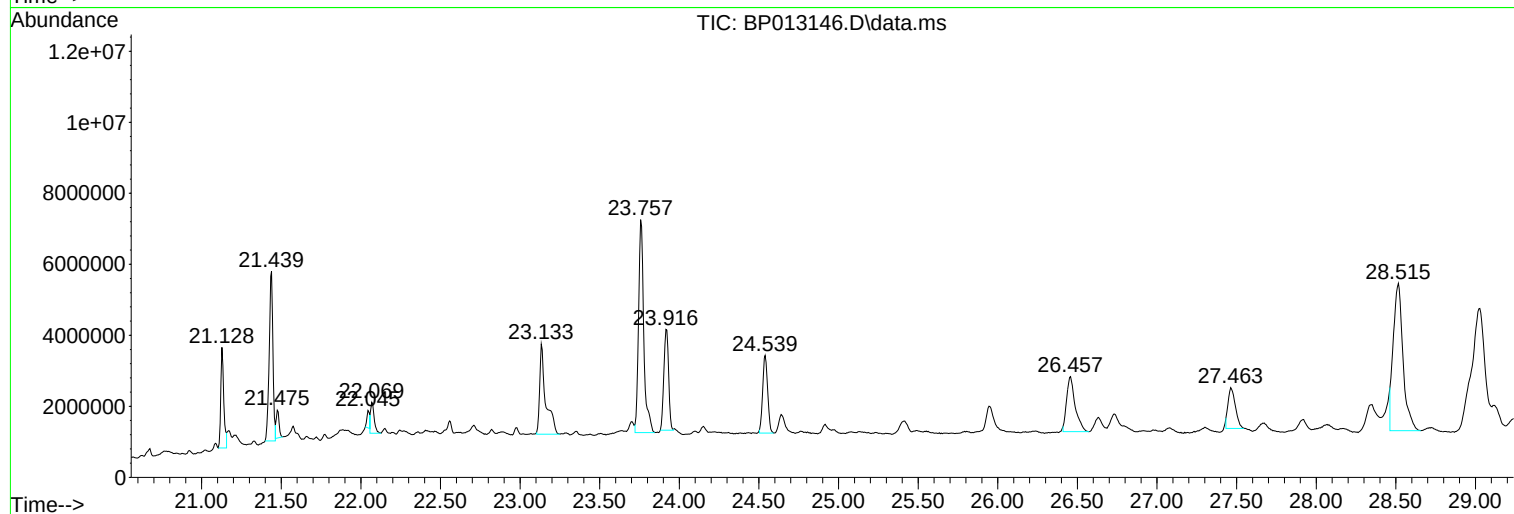
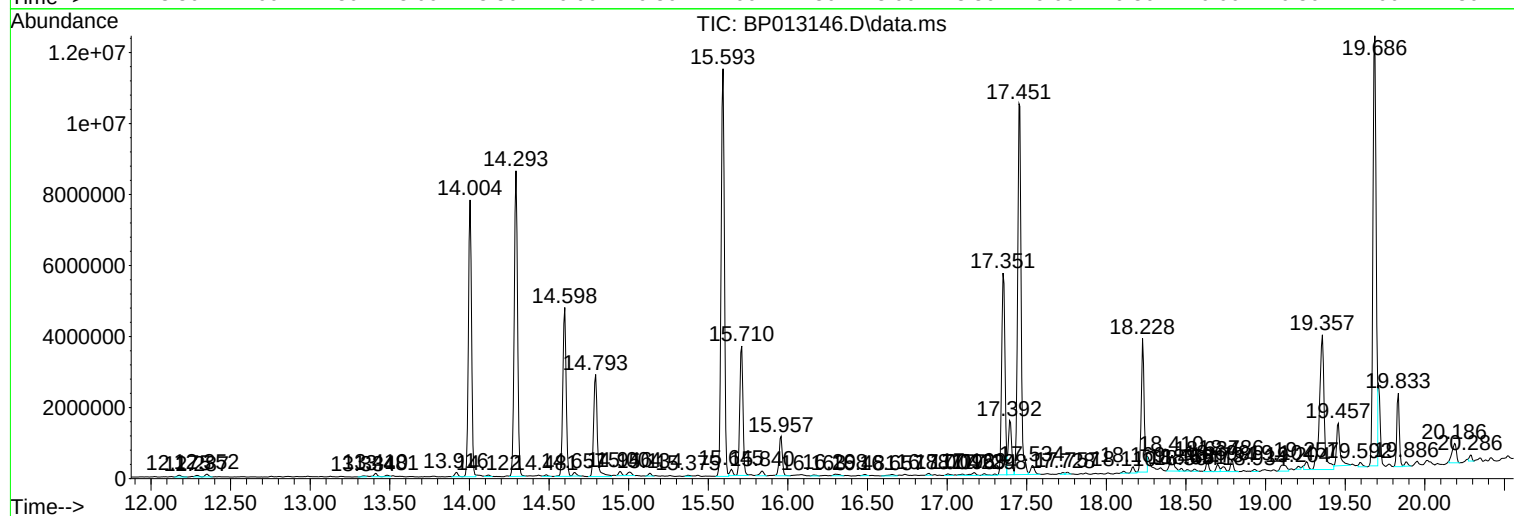
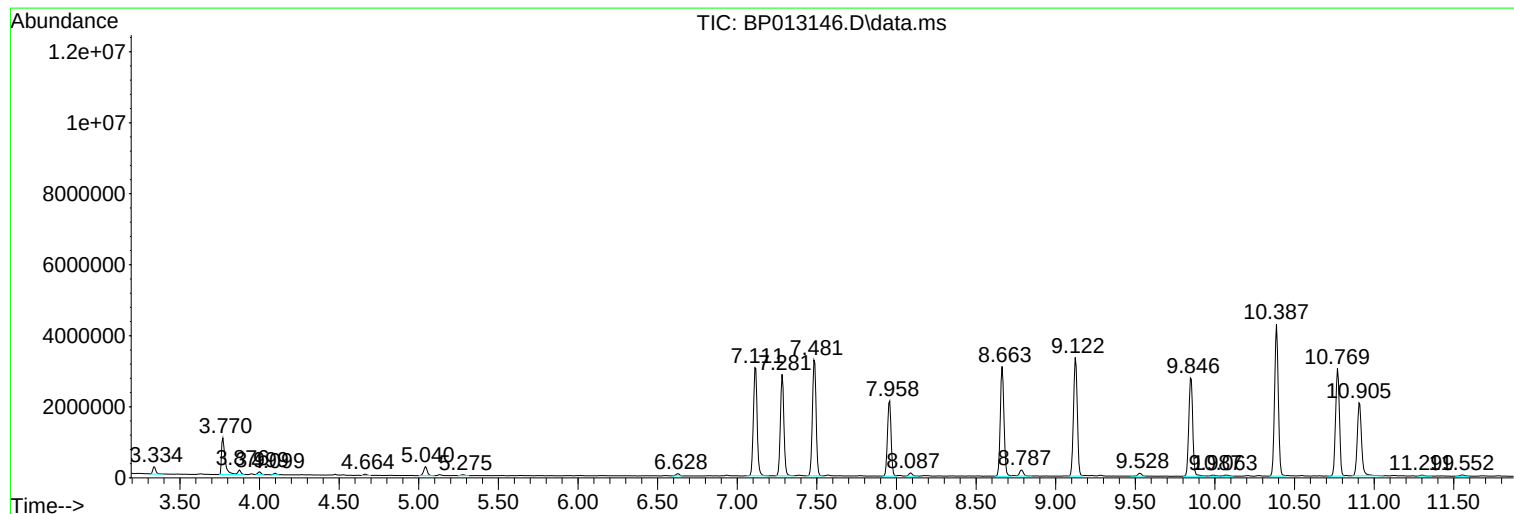
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
Quant Title : SVOA CALIBRATION

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



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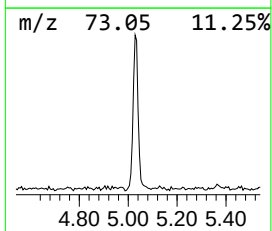
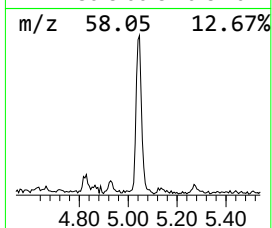
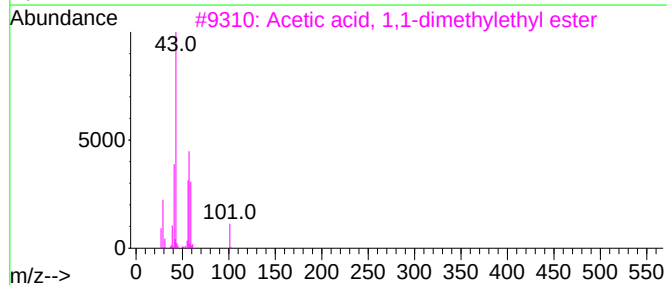
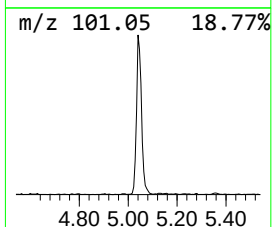
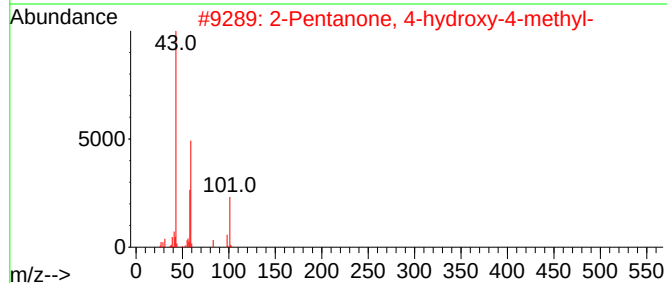
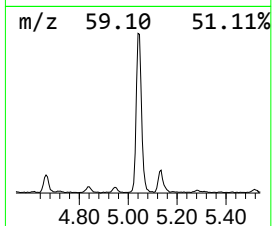
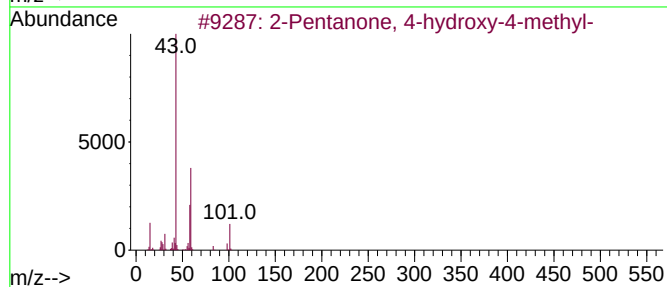
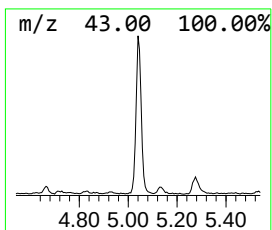
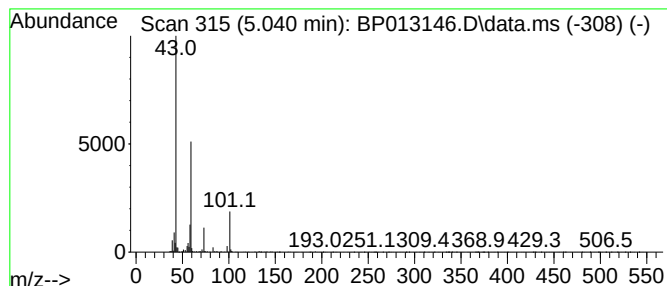
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.040	2.48 ng/ul	416077	1,4-Dichlorobenzene-d4	7.958

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	59
2			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50
3			Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	37
4			1-(1-Methoxypropan-2-yloxy)propa...	260	C14H28O4	1000367-13-1	37
5			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	36



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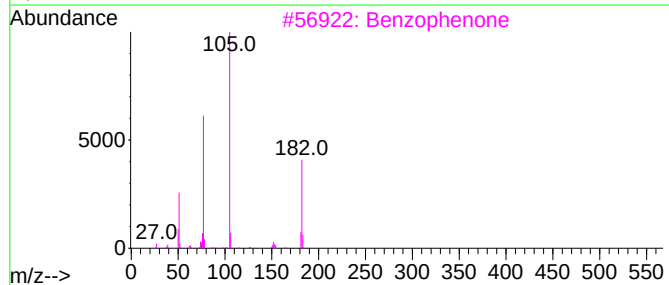
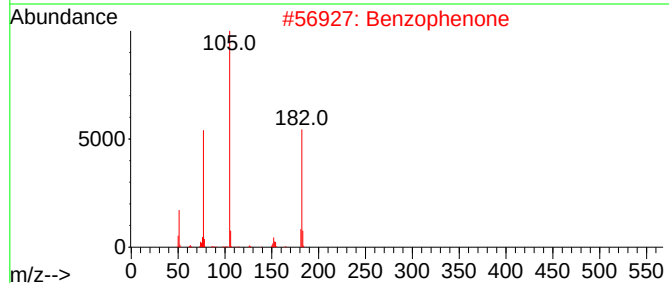
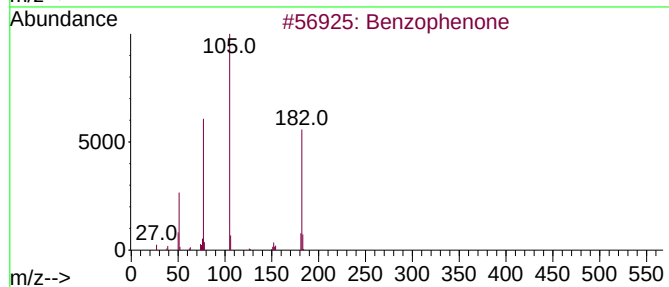
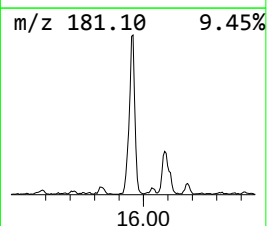
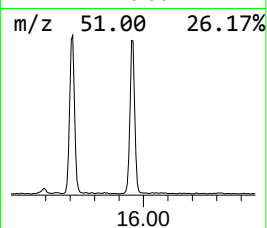
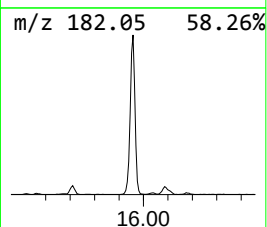
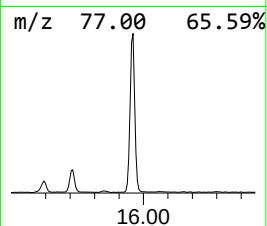
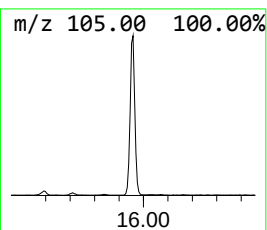
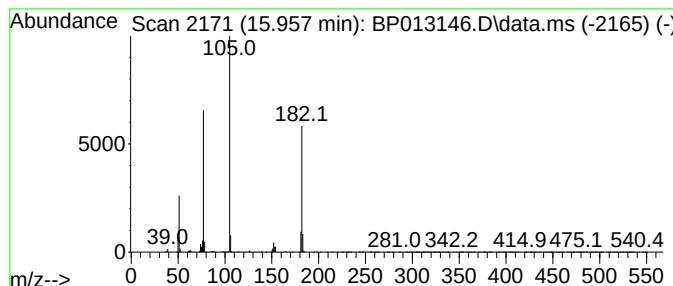
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 Benzophenone Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.957	4.32 ng/ul	1501860	Acenaphthene-d10	14.599

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzophenone	182	C13H10O	000119-61-9	96
2			Benzophenone	182	C13H10O	000119-61-9	95
3			Benzophenone	182	C13H10O	000119-61-9	94
4			Benzophenone	182	C13H10O	000119-61-9	91
5			Benzophenone	182	C13H10O	000119-61-9	90



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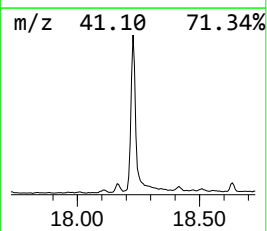
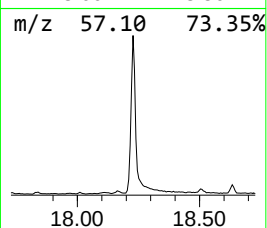
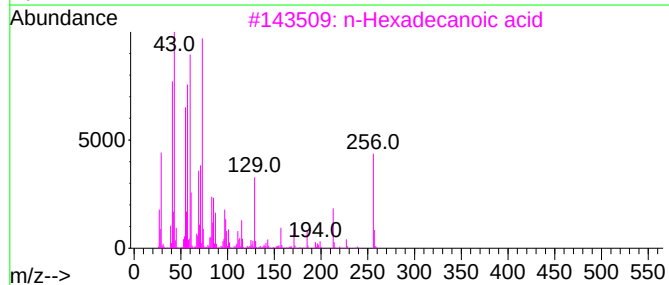
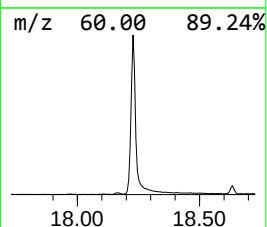
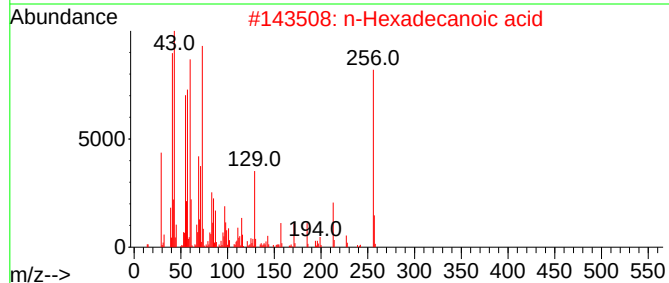
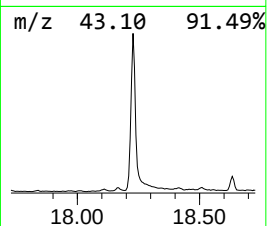
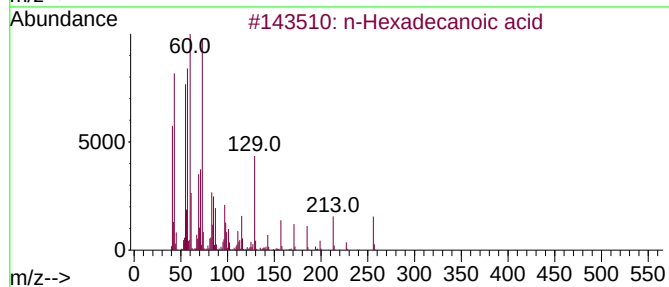
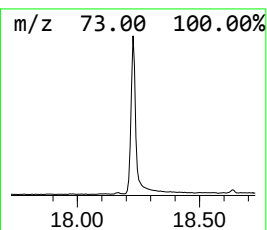
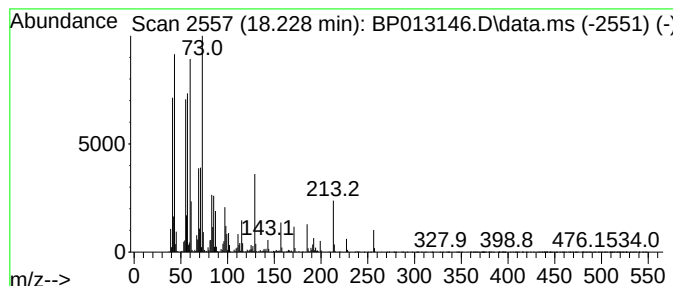
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 n-Hexadecanoic acid Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.228	12.12 ng/ul	4979140	Phenanthrene-d10	17.351

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	95
5			Tridecanoic acid	214	C13H26O2	000638-53-9	95



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121922\
Data File : BP013146.D
Acq On : 20 Dec 2022 02:18
Operator : CG/JU
Sample : N6006-06
Misc :
ALS Vial : 25 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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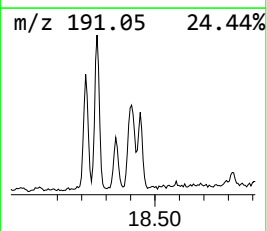
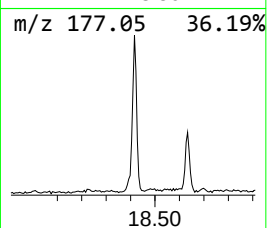
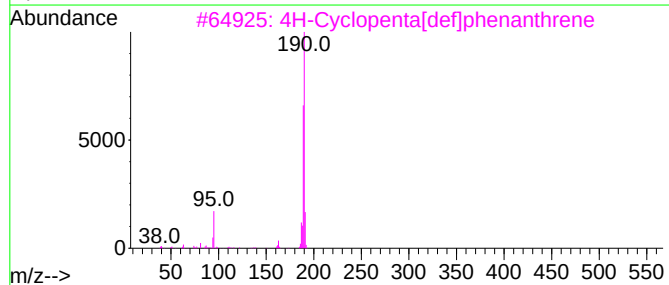
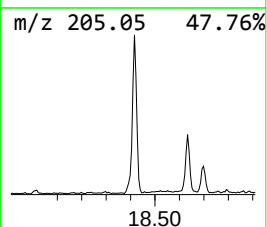
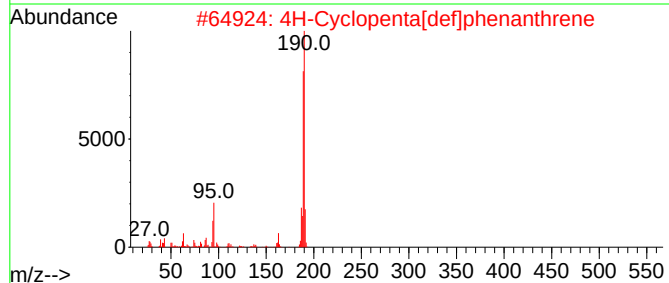
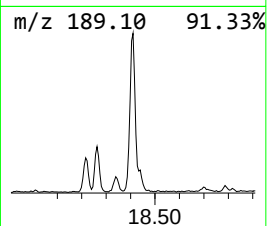
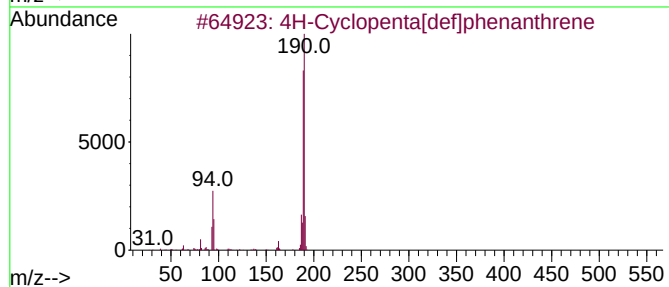
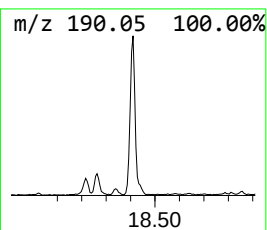
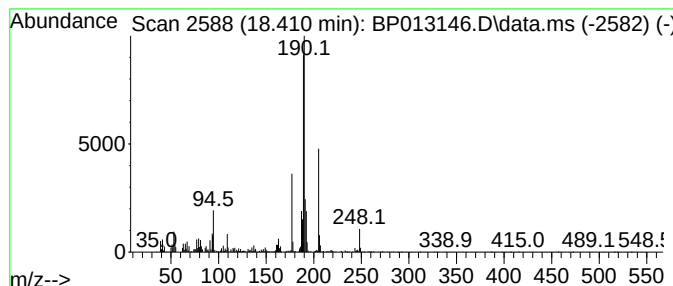
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 4H-Cyclopenta[def]phenanthrene Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.410	2.01 ng/ul	827148	Phenanthrene-d10	17.351

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	64
2			4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	64
3			4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	52
4			2,3,5,6-Tetramethylterephthalald...	190	C12H14O2	007072-01-7	49
5			6H-Cyclobuta[jk]phenanthrene	190	C15H10	083469-43-6	40



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121922\
Data File : BP013146.D
Acq On : 20 Dec 2022 02:18
Operator : CG/JU
Sample : N6006-06
Misc :
ALS Vial : 25 Sample Multiplier: 1

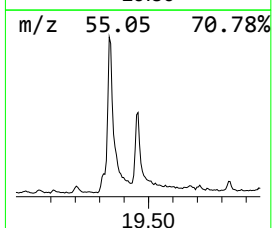
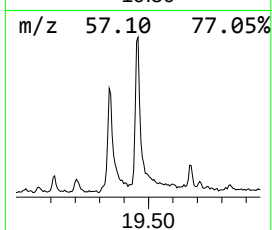
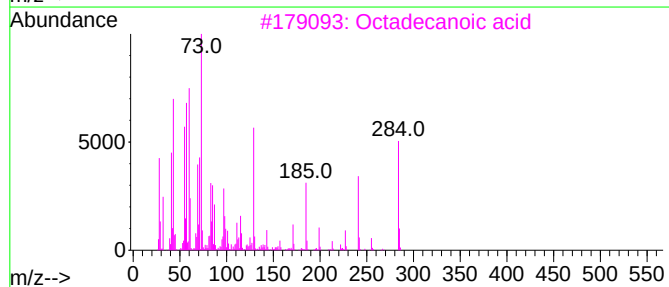
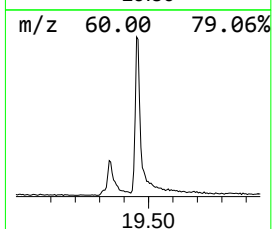
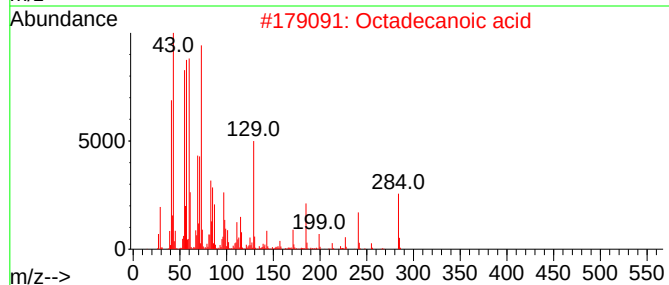
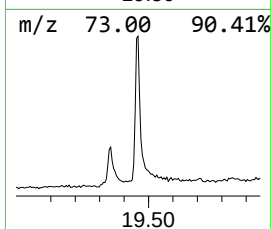
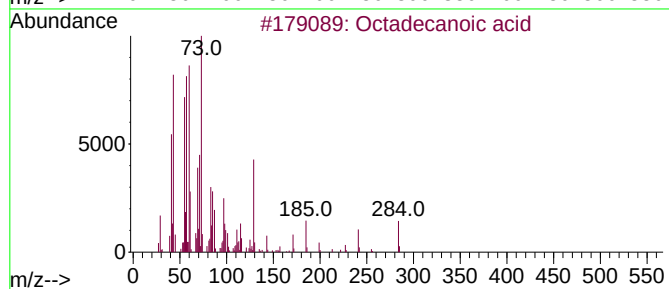
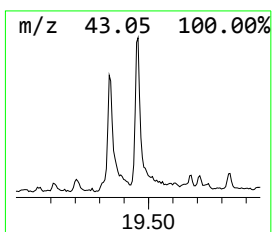
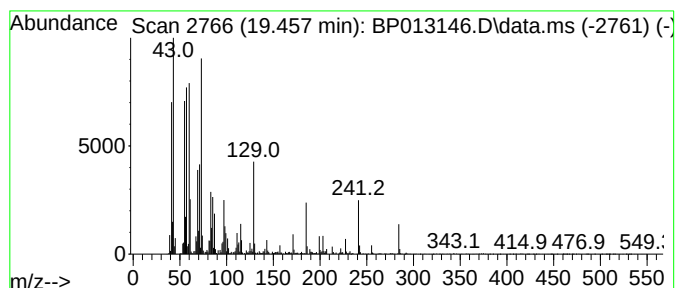
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 Octadecanoic acid Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.	
19.457	4.52 ng/ul	1780260	Chrysene-d12	21.439	
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	97
4	Octadecanoic acid	284	C18H36O2	000057-11-4	90
5	Octadecanoic acid	284	C18H36O2	000057-11-4	86



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121922\
Data File : BP013146.D
Acq On : 20 Dec 2022 02:18
Operator : CG/JU
Sample : N6006-06
Misc :
ALS Vial : 25 Sample Multiplier: 1

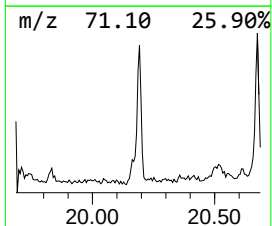
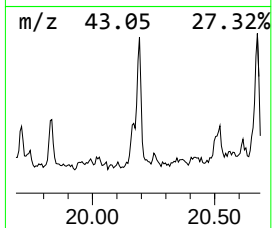
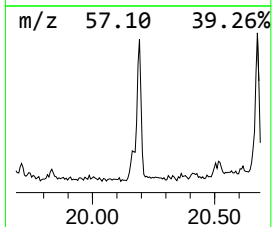
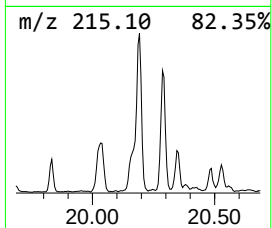
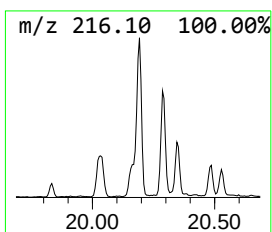
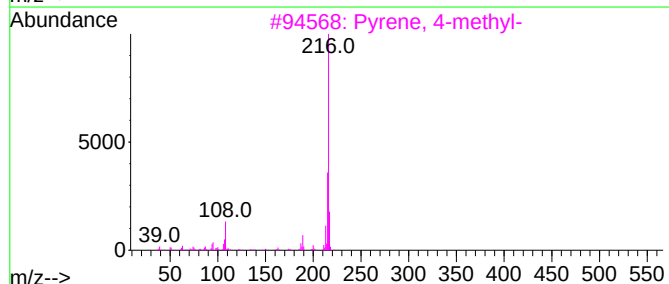
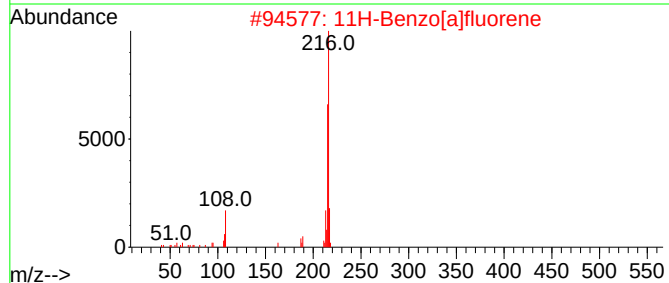
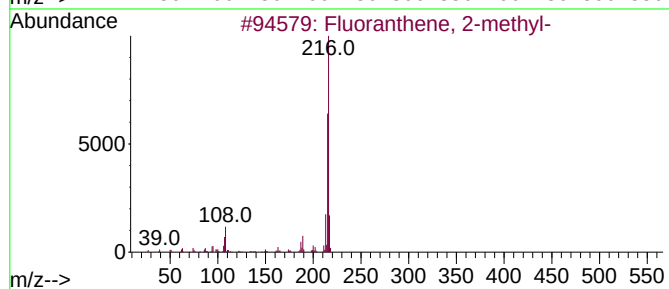
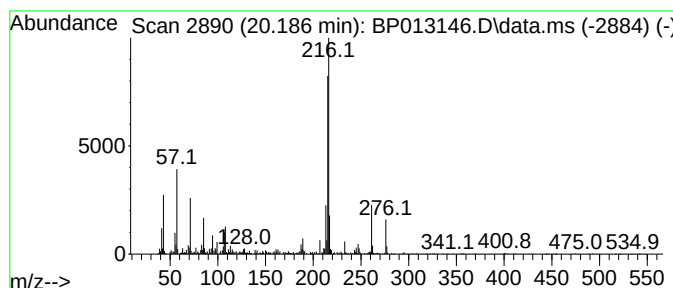
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
Quant Title : SVOA CALIBRATION

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

Peak Number 7 Fluoranthene, 2-methyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.	
20.186	2.65 ng/ul	1044250	Chrysene-d12	21.439	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	96
2	11H-Benzo[a]fluorene	216	C17H12	000238-84-6	90
3	Pyrene, 4-methyl-	216	C17H12	003353-12-6	89
4	11H-Benzo[b]fluorene	216	C17H12	000243-17-4	89
5	11H-Benzo[a]fluorene	216	C17H12	000238-84-6	81



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121922\
Data File : BP013146.D
Acq On : 20 Dec 2022 02:18
Operator : CG/JU
Sample : N6006-06
Misc :
ALS Vial : 25 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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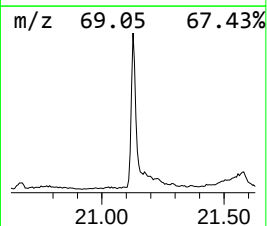
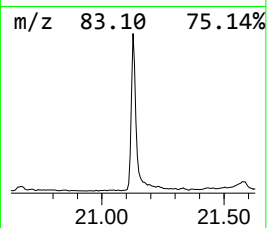
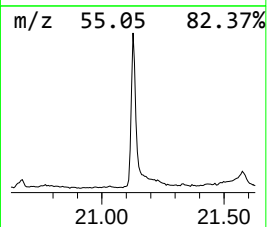
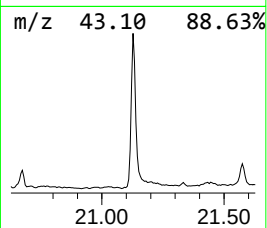
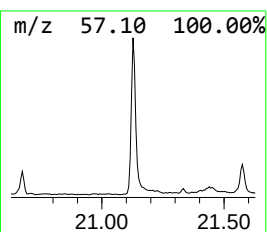
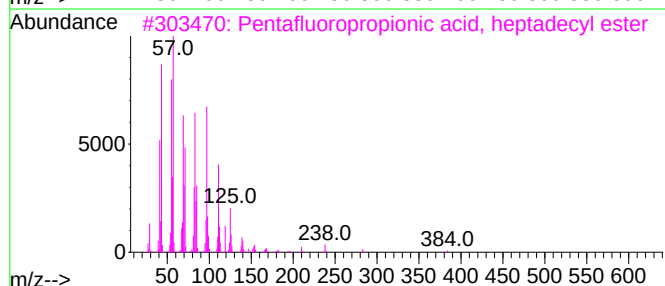
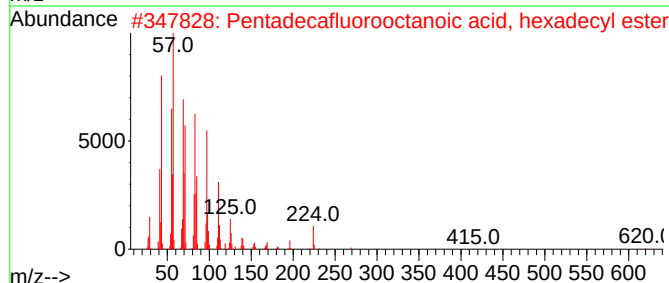
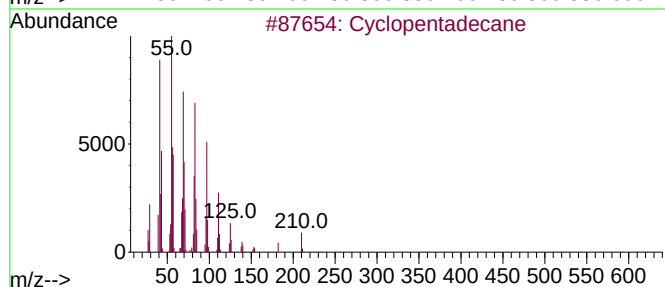
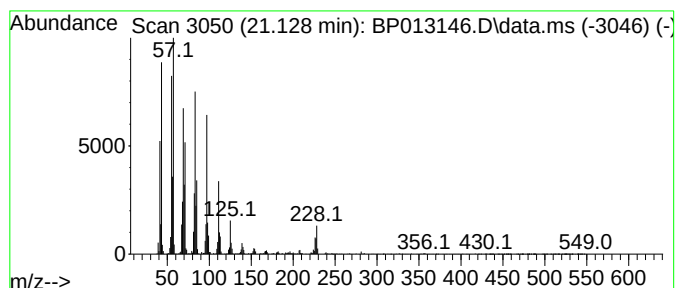
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 (DEL) Alkane: Cyclic21.128 Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.128	9.54 ng/ul	3758130	Chrysene-d12	21.439

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclopentadecane	210	C15H30	000295-48-7	97
2			Pentadecafluorooctanoic acid, he...	638	C24H33F15O2	1000406-04-6	94
3			Pentafluoropropionic acid, hepta...	402	C20H35F5O2	959218-78-1	94
4			Trifluoroacetoxy hexadecane	338	C18H33F3O2	006222-03-3	94
5			Cyclotetracosane	336	C24H48	000297-03-0	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121922\
Data File : BP013146.D
Acq On : 20 Dec 2022 02:18
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Sample : N6006-06
Misc :
ALS Vial : 25 Sample Multiplier: 1

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ClientSampleId :
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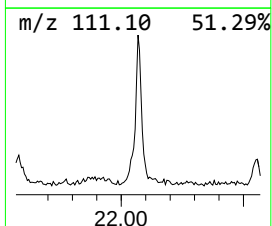
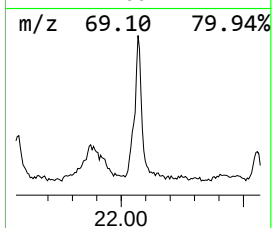
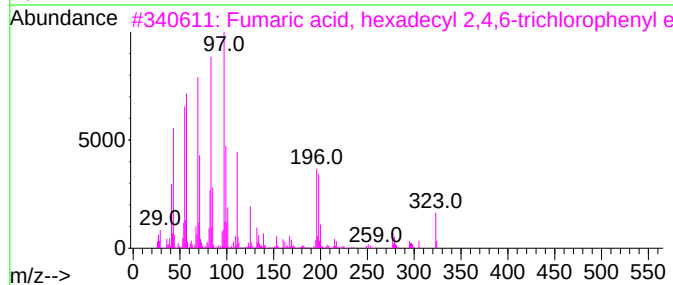
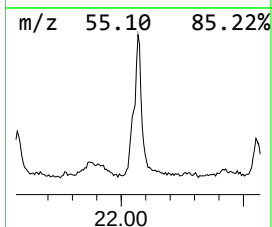
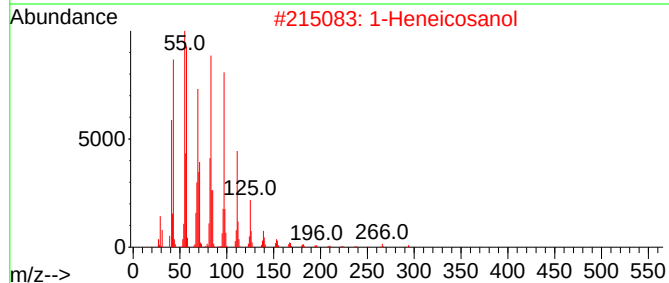
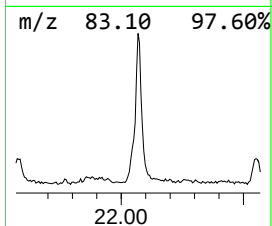
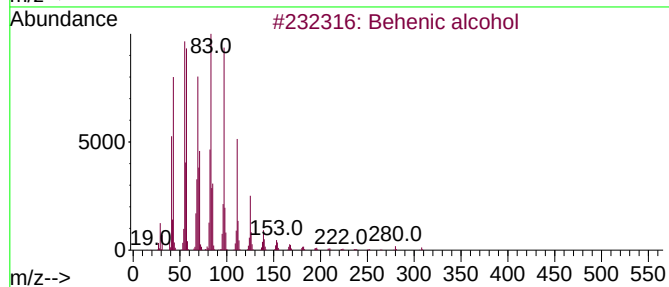
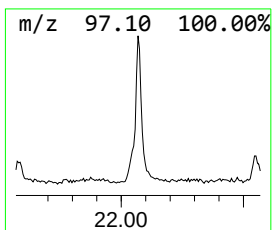
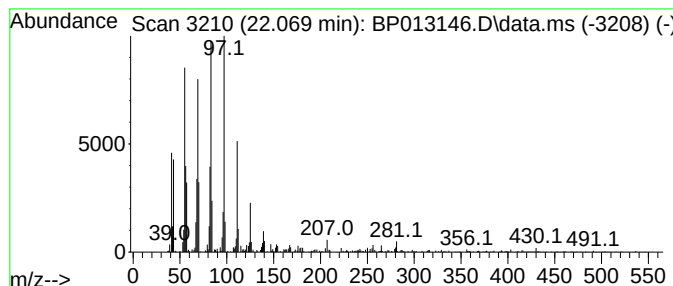
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 Behenic alcohol Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.069	3.25 ng/ul	1279780	Chrysene-d12	21.439

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Behenic alcohol	326	C22H46O	000661-19-8	94
2			1-Heneicosanol	312	C21H44O	015594-90-8	83
3			Fumaric acid, hexadecyl 2,4,6-tr...	518	C26H37Cl3O4	1000348-27-9	80
4			1-Docosanol, methyl ether	340	C23H48O	1000333-91-9	74
5			Dodecane, 1-cyclopentyl-4-(3-cyc...	348	C25H48	007225-68-5	74



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121922\
Data File : BP013146.D
Acq On : 20 Dec 2022 02:18
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Sample : N6006-06
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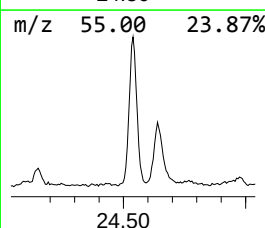
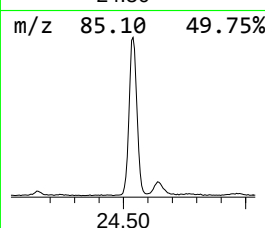
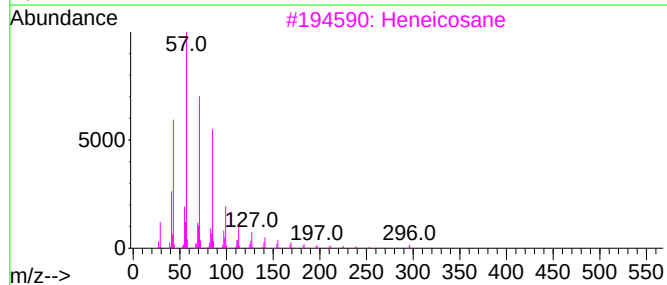
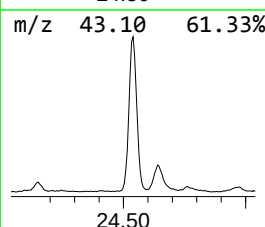
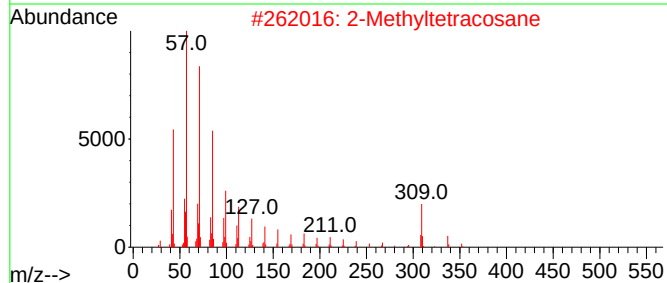
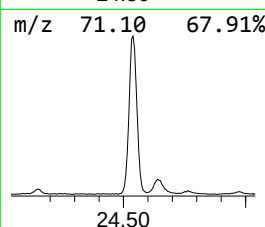
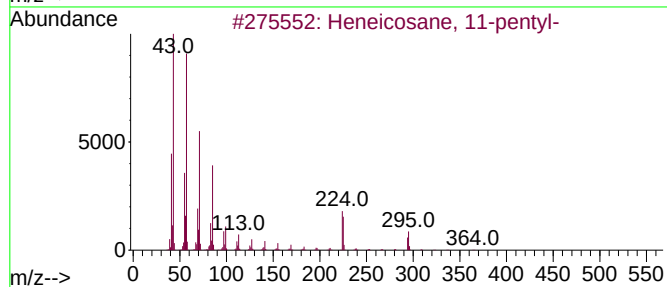
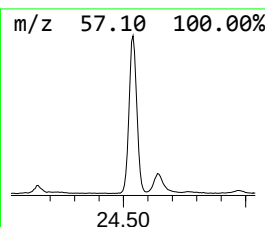
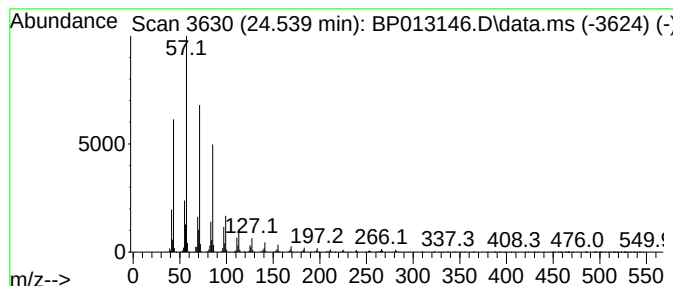
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 (DEL) Alkane: Straight-Chai... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.539	16.19 ng/ul	4730340	Perylene-d12	23.922

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Heneicosane, 11-pentyl-	366	C26H54	014739-72-1	93
2		2-Methyltetracosane	352	C25H52	001560-78-7	93
3		Heneicosane	296	C21H44	000629-94-7	91
4		Octacosane, 2-methyl-	408	C29H60	001560-98-1	91
5		Triacontane	422	C30H62	000638-68-6	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121922\
Data File : BP013146.D
Acq On : 20 Dec 2022 02:18
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Sample : N6006-06
Misc :
ALS Vial : 25 Sample Multiplier: 1

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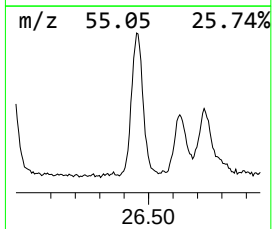
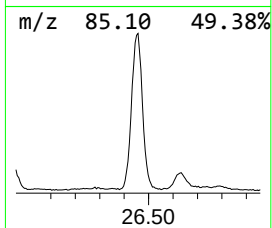
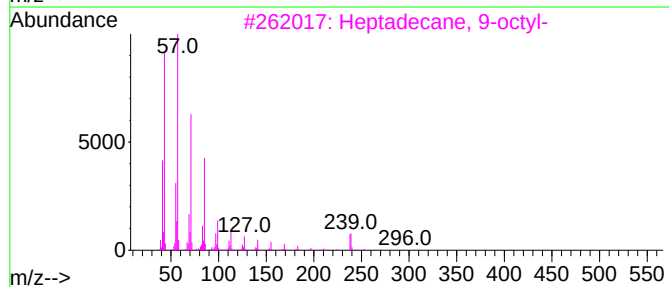
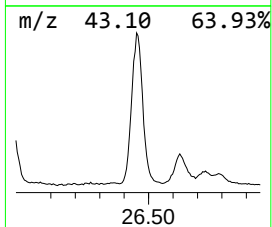
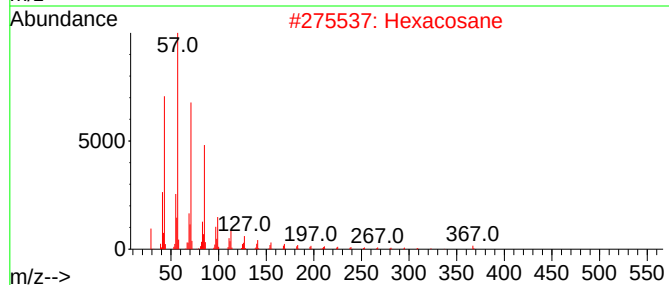
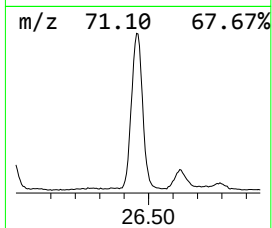
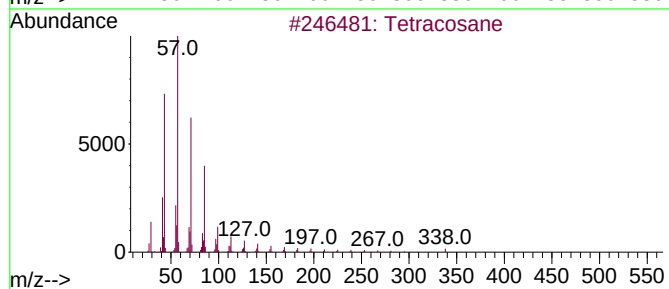
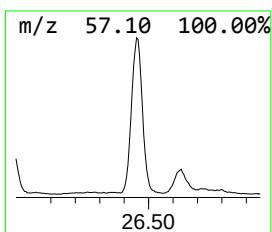
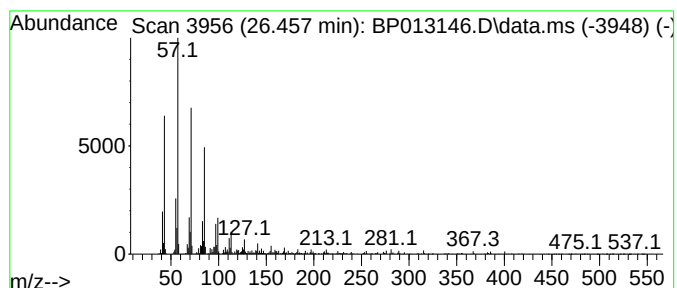
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Quant Title : SVOA CALIBRATION

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

Peak Number 11 (DEL) Alkane: Straight-Chai... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
26.457	19.68 ng/ul	5751420	Perylene-d12	23.922

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tetracosane	338	C24H50	000646-31-1	95
2			Hexacosane	366	C26H54	000630-01-3	94
3			Heptadecane, 9-octyl-	352	C25H52	007225-64-1	93
4			2-Methylheptacosane	394	C28H58	001561-00-8	93
5			Eicosane	282	C20H42	000112-95-8	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121922\
Data File : BP013146.D
Acq On : 20 Dec 2022 02:18
Operator : CG/JU
Sample : N6006-06
Misc :
ALS Vial : 25 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DBXV9

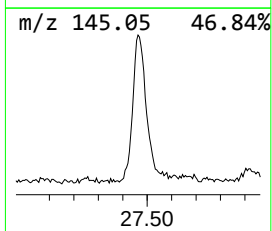
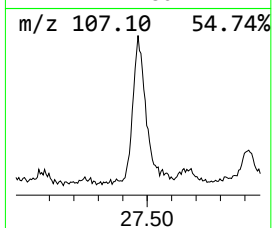
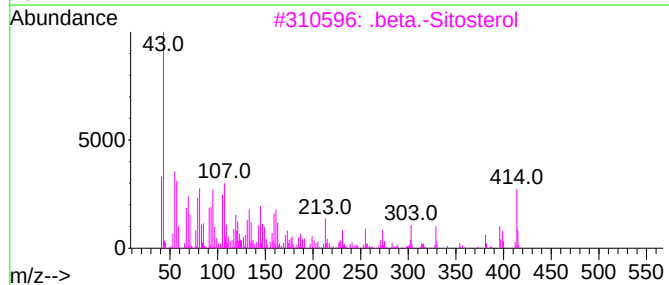
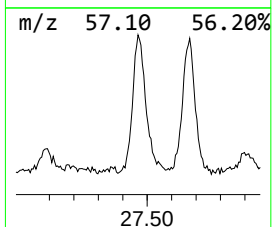
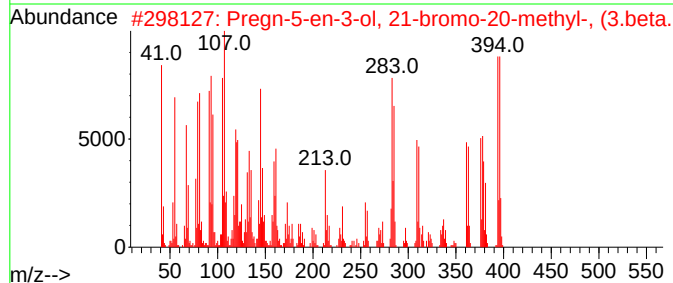
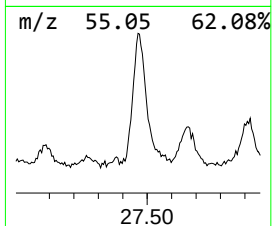
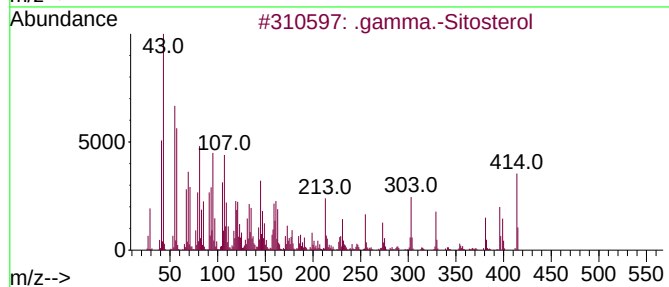
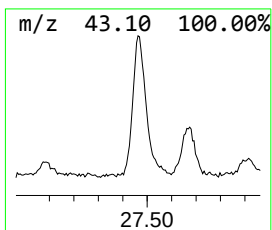
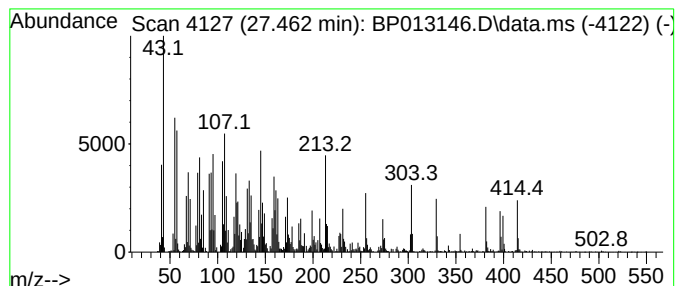
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Quant Title : SVOA CALIBRATION

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

Peak Number 12 .gamma.-Sitosterol Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
27.462	12.38 ng/ul	3616990	Perylene-d12	23.922

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	.	gamma.-Sitosterol	414	C29H50O	000083-47-6	99	
2	Pregn-5-en-3-ol, 21-bromo-20-met...	394	C22H35BrO	055103-80-5	93		
3	.beta.-Sitosterol	414	C29H50O	000083-46-5	92		
4	.beta.-Sitosterol	414	C29H50O	000083-46-5	83		
5	Ergost-5-en-3-ol, (3.beta.)-	400	C28H48O	004651-51-8	49		



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121922\
Data File : BP013146.D
Acq On : 20 Dec 2022 02:18
Operator : CG/JU
Sample : N6006-06
Misc :
ALS Vial : 25 Sample Multiplier: 1

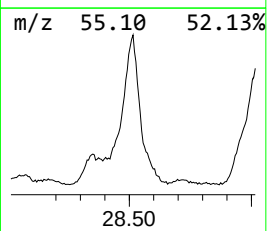
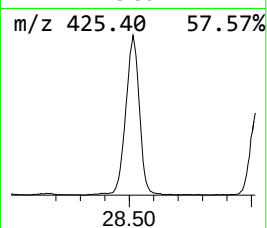
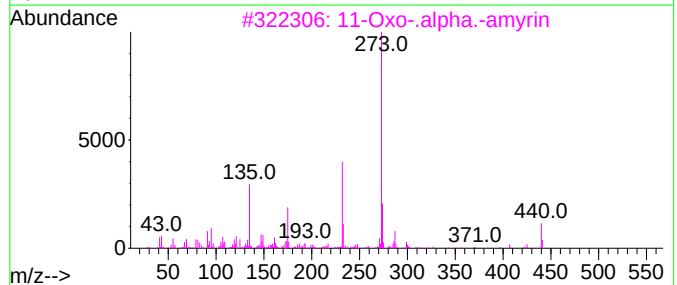
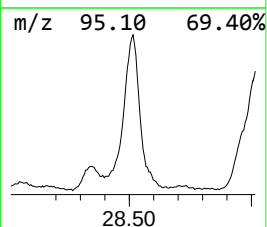
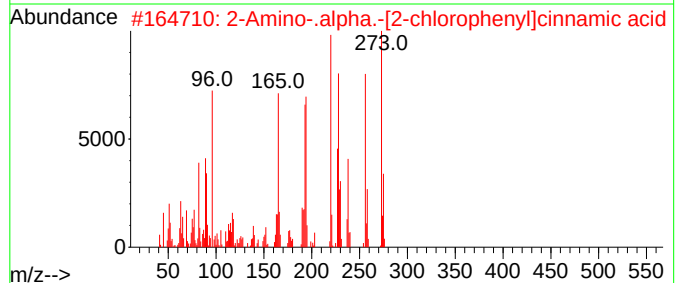
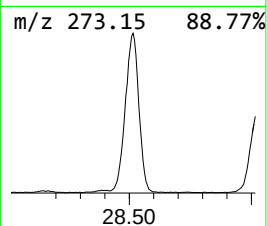
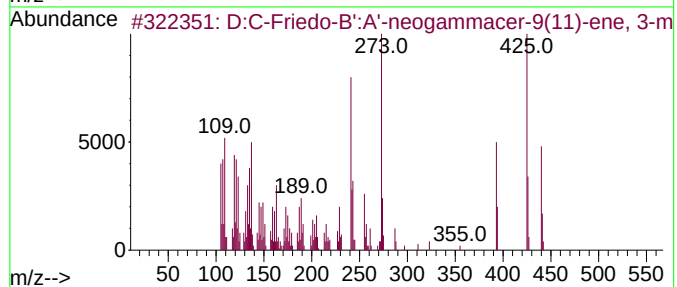
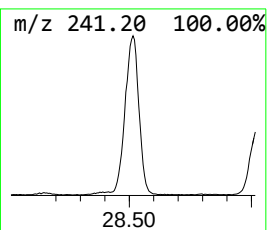
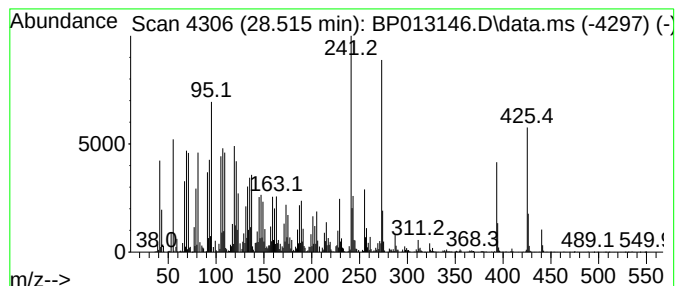
Instrument :
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
Quant Title : SVOA CALIBRATION

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

Peak Number 13 D:C-Friedo-B':A'-neogammace... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.		
28.515	60.41 ng/ul	17652000	Perylene-d12	23.922		
Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		D:C-Friedo-B':A'-neogammacer-9(1...	440	C31H52O	004555-56-0	78
2		2-Amino-.alpha.-[2-chlorophenyl]...	273	C15H12ClNO2	1000254-08-5	59
3		11-Oxo-.alpha.-amyrin	440	C30H48O2	002118-90-3	45
4		2-(1H-Indol-3-ylmethylene)-1H-in...	273	C18H11NO2	031060-64-7	25
5		Pyrido[3,4-b]benzothiophene-4-ca...	273	C13H8FN3OS	1000265-50-3	15



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121922\
Data File : BP013146.D
Acq On : 20 Dec 2022 02:18
Operator : CG/JU
Sample : N6006-06
Misc :
ALS Vial : 25 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DBXV9

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.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
2-Pentanone, 4-...	5.040	2.5	ng/ul	416077	1	7.958	3351230	20.0
Benzophenone	15.957	4.3	ng/ul	1501860	3	14.599	6951190	20.0
n-Hexadecanoic ...	18.228	12.1	ng/ul	4979140	4	17.351	8213390	20.0
4H-Cyclopenta[d...]	18.410	2.0	ng/ul	827148	4	17.351	8213390	20.0
Octadecanoic acid	19.457	4.5	ng/ul	1780260	5	21.439	7882270	20.0
(1H)Pyrrole, 3,...	19.834	5.9	ng/ul	2326480	5	21.439	7882270	20.0
Fluoranthene, 2...	20.186	2.6	ng/ul	1044250	5	21.439	7882270	20.0
(DEL) Alkane: C...	21.128	9.5	ng/ul	3758130	5	21.439	7882270	20.0
Behenic alcohol	22.069	3.3	ng/ul	1279780	5	21.439	7882270	20.0
(DEL) Alkane: S...	24.539	16.2	ng/ul	4730340	6	23.922	5844260	20.0
(DEL) Alkane: S...	26.457	19.7	ng/ul	5751420	6	23.922	5844260	20.0
.gamma.-Sitosterol	27.462	12.4	ng/ul	3616990	6	23.922	5844260	20.0
D:C-Friedo-B':A...	28.515	60.4	ng/ul	17652000	6	23.922	5844260	20.0