

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122222\
 Data File : BP013243.D
 Acq On : 22 Dec 2022 13:22
 Operator : CG/JU
 Sample : N6015-02
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 DBXT6

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: BP013243.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	21	25	33	rVB	146466	194075	0.93%	0.085%
2	3.764	95	98	114	rBV	1596614	2430235	11.66%	1.063%
3	4.093	150	154	160	rVB	53417	79322	0.38%	0.035%
4	4.464	215	217	224	rVB5	29976	39447	0.19%	0.017%
5	5.040	307	315	321	rBV3	35069	75090	0.36%	0.033%
6	5.264	349	353	362	rBV	44842	83114	0.40%	0.036%
7	6.122	491	499	504	rBV	17320	41257	0.20%	0.018%
8	6.617	576	583	589	rVV2	50132	83198	0.40%	0.036%
9	7.105	658	666	679	rBV	2470329	3825734	18.36%	1.674%
10	7.275	689	695	702	rBV	1946053	2996412	14.38%	1.311%
11	7.475	722	729	738	rBV	2532648	3888996	18.66%	1.701%
12	7.946	800	809	816	rBV	2472998	3796197	18.22%	1.661%
13	8.658	922	930	944	rBV	2612792	4229415	20.30%	1.850%
14	9.116	1000	1008	1015	rBV	2352588	3764503	18.07%	1.647%
15	9.269	1028	1034	1040	rBV	74915	115940	0.56%	0.051%
16	9.516	1070	1076	1082	rVB	71842	113522	0.54%	0.050%
17	9.840	1122	1131	1144	rBV	1966066	3232310	15.51%	1.414%
18	10.375	1214	1222	1237	rBV	3103002	5106566	24.51%	2.234%
19	10.757	1278	1287	1295	rBV	3283639	5602685	26.89%	2.451%
20	10.899	1304	1311	1329	rVB	1559832	2758168	13.24%	1.207%
21	11.293	1372	1378	1385	rBV7	18046	36139	0.17%	0.016%
22	12.010	1495	1500	1505	rBV6	17462	32238	0.15%	0.014%
23	12.075	1505	1511	1514	rBV2	18821	31466	0.15%	0.014%
24	12.169	1522	1527	1535	rVB	44106	67626	0.32%	0.030%
25	12.340	1552	1556	1560	rBV	46818	62987	0.30%	0.028%
26	12.857	1638	1644	1649	rBV8	27906	52618	0.25%	0.023%
27	13.328	1718	1724	1726	rBV7	26087	39839	0.19%	0.017%
28	13.404	1730	1737	1741	rVV	90867	151058	0.72%	0.066%
29	13.457	1741	1746	1751	rVV9	19876	44619	0.21%	0.020%
30	13.504	1751	1754	1759	rVV3	32460	62108	0.30%	0.027%
31	13.575	1762	1766	1771	rVB	58769	86078	0.41%	0.038%
32	13.910	1818	1823	1829	rVV7	40022	72128	0.35%	0.032%
33	13.993	1829	1837	1845	rVV	4795667	6951182	33.36%	3.041%
34	14.057	1845	1848	1853	rVV3	45637	76633	0.37%	0.034%
35	14.110	1853	1857	1861	rVV6	33413	56164	0.27%	0.025%
36	14.281	1879	1886	1898	rBV	5850240	8527440	40.93%	3.731%

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Integrator: RTE

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

37	14.475	1915	1919	1926	rVB	64365	97798	0.47%	0.043%
38	14.587	1929	1938	1945	rBV	5044969	7447465	35.74%	3.258%
39	14.651	1945	1949	1955	rVB6	35117	52037	0.25%	0.023%
40	14.787	1965	1972	1978	rBV	1532013	2477874	11.89%	1.084%
41	14.940	1995	1998	2002	rVB3	55667	65831	0.32%	0.029%
42	14.993	2002	2007	2012	rVV3	65281	100304	0.48%	0.044%
43	15.040	2012	2015	2025	rVB3	24530	46854	0.22%	0.020%
44	15.422	2077	2080	2084	rVB3	31414	47412	0.23%	0.021%
45	15.498	2088	2093	2097	rVB5	36130	50728	0.24%	0.022%
46	15.581	2101	2107	2117	rVV	7418852	10539404	50.58%	4.611%
47	15.698	2121	2127	2137	rBV	1971426	2618091	12.56%	1.145%
48	15.822	2144	2148	2155	rVB4	39794	66901	0.32%	0.029%
49	15.945	2159	2169	2181	rBV	2607409	3557126	17.07%	1.556%
50	16.151	2197	2204	2210	rVB	32610	72628	0.35%	0.032%
51	16.216	2211	2215	2221	rVV5	47980	70590	0.34%	0.031%
52	16.292	2224	2228	2236	rVV4	55490	123400	0.59%	0.054%
53	16.363	2236	2240	2245	rVV3	22955	44840	0.22%	0.020%
54	16.434	2248	2252	2254	rBV5	37174	51550	0.25%	0.023%
55	16.475	2256	2259	2263	rVV	58900	92749	0.45%	0.041%
56	16.516	2263	2266	2274	rVB2	85890	123014	0.59%	0.054%
57	16.728	2298	2302	2312	rBV3	138741	222360	1.07%	0.097%
58	16.828	2315	2319	2325	rVB3	29672	61130	0.29%	0.027%
59	17.092	2360	2364	2367	rBV4	43187	63527	0.30%	0.028%
60	17.228	2383	2387	2391	rBV	144923	187180	0.90%	0.082%
61	17.292	2395	2398	2401	rBV3	69957	81923	0.39%	0.036%
62	17.345	2401	2407	2417	rVV	5664800	8512619	40.85%	3.724%
63	17.445	2417	2424	2431	rVB	6971219	9926839	47.64%	4.343%
64	17.557	2440	2443	2449	rVB7	51447	64742	0.31%	0.028%
65	17.622	2451	2454	2459	rVB6	54785	80614	0.39%	0.035%
66	17.698	2464	2467	2482	rVB6	102792	210280	1.01%	0.092%
67	17.828	2486	2489	2492	rBV2	46845	52755	0.25%	0.023%
68	17.869	2492	2496	2499	rBV5	28695	40884	0.20%	0.018%
69	18.004	2515	2519	2522	rBV4	68913	79712	0.38%	0.035%
70	18.045	2522	2526	2530	rBV3	65071	95435	0.46%	0.042%
71	18.104	2530	2536	2540	rBV	149218	248475	1.19%	0.109%
72	18.157	2541	2545	2550	rBV2	225944	327582	1.57%	0.143%
73	18.222	2550	2556	2569	rBV	3300305	4670052	22.41%	2.043%
74	18.410	2581	2588	2593	rVB	701447	943869	4.53%	0.413%
75	18.463	2593	2597	2601	rBV2	310427	439410	2.11%	0.192%
76	18.628	2621	2625	2630	rBV	373571	447996	2.15%	0.196%

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Integrator: RTE

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 Title : SVOA CALIBRATION

77	18.716	2638	2640	2647	rVB	113897	161963	0.78%	0.071%
78	18.781	2647	2651	2658	rVB	247997	337586	1.62%	0.148%
79	19.122	2703	2709	2713	rBV2	539296	740258	3.55%	0.324%
80	19.251	2728	2731	2735	rVB3	117937	142677	0.68%	0.062%
81	19.310	2737	2741	2742	rBV	291518	332068	1.59%	0.145%
82	19.333	2742	2745	2747	rBV2	331743	395609	1.90%	0.173%
83	19.457	2760	2766	2786	rBV	12744625	17621922	84.57%	7.709%
84	19.675	2798	2803	2816	rVB	7200037	10117296	48.56%	4.426%
85	19.828	2824	2829	2840	rBV	3688333	4341463	20.84%	1.899%
86	20.169	2883	2887	2894	rVB4	674202	1031340	4.95%	0.451%
87	20.498	2940	2943	2949	rVB4	180820	301566	1.45%	0.132%
88	21.128	3045	3050	3064	rBV2	2337751	5876618	28.20%	2.571%
89	21.433	3097	3102	3107	rBV	4911399	6925150	33.24%	3.030%
90	22.039	3202	3205	3207	rBV	402804	502756	2.41%	0.220%
91	22.069	3207	3210	3220	rVB2	743829	1194547	5.73%	0.523%
92	23.122	3385	3389	3395	rBV	1164327	1793615	8.61%	0.785%
93	23.745	3489	3495	3510	rVB2	4388163	9503682	45.61%	4.158%
94	23.910	3516	3523	3536	rVB2	3167987	6693721	32.13%	2.928%
95	24.527	3622	3628	3636	rVB	1957496	4188066	20.10%	1.832%
96	24.627	3640	3645	3660	rVB2	1479061	4065949	19.51%	1.779%
97	25.198	3737	3742	3752	rVB	1021497	2428443	11.65%	1.062%
98	26.439	3944	3953	3972	rVB	2713863	9034818	43.36%	3.952%
99	27.451	4117	4125	4137	rBV2	1945729	6707264	32.19%	2.934%
100	29.004	4378	4389	4416	rVB9	3980577	20836386	100.00%	9.115%

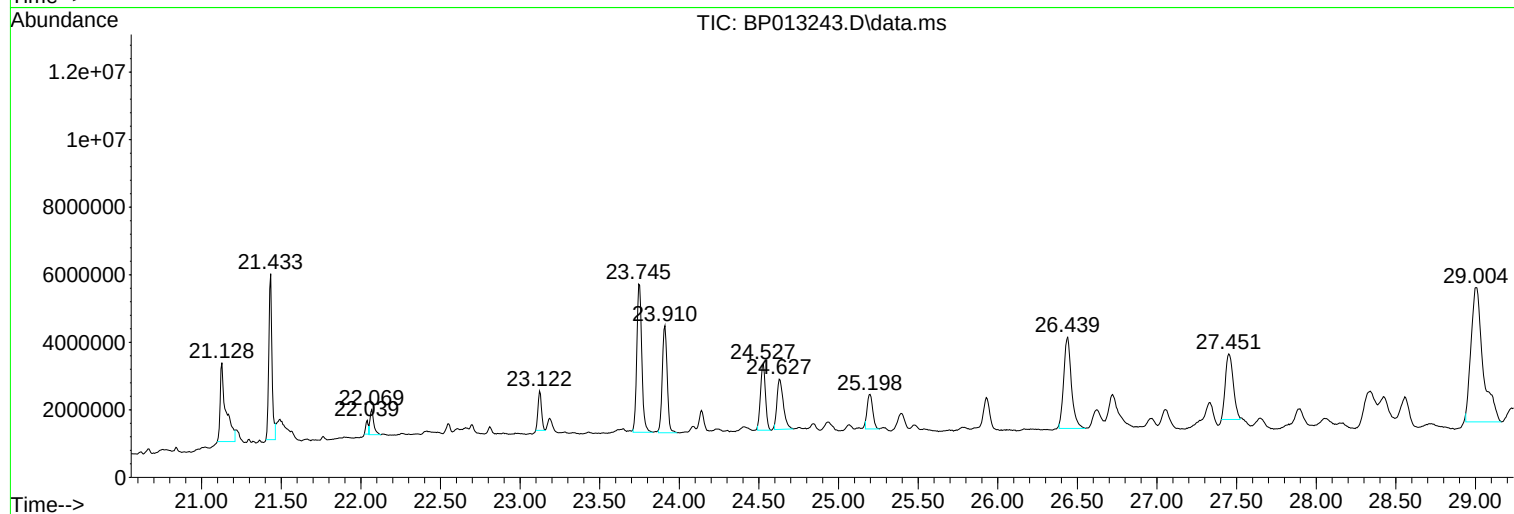
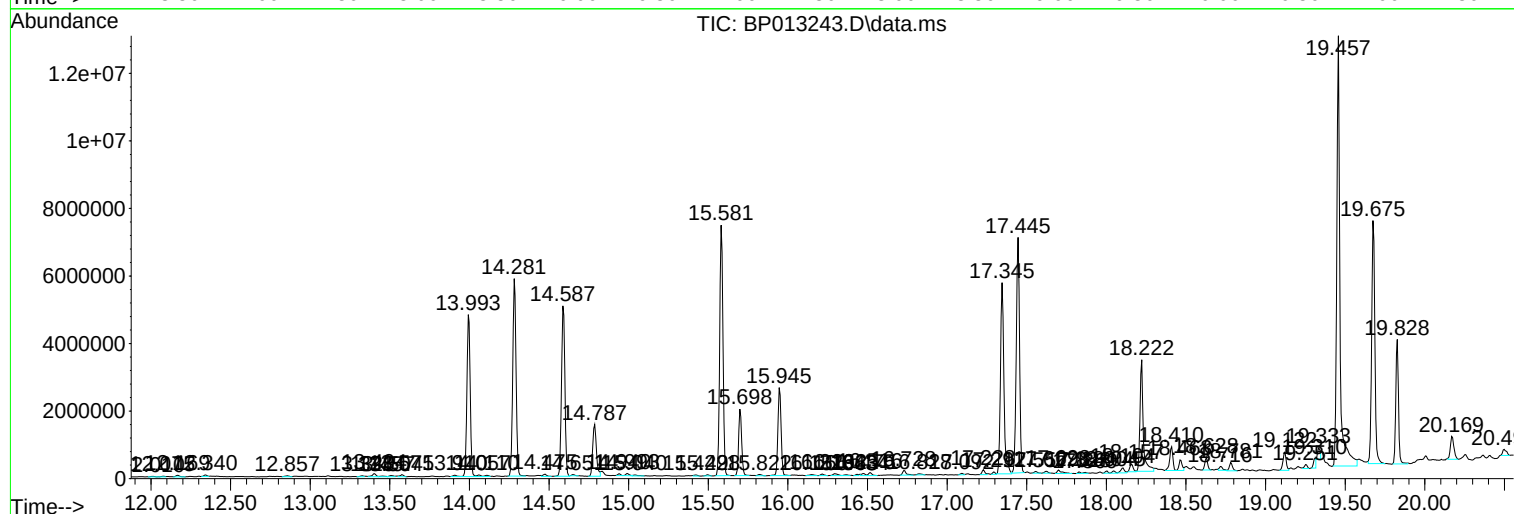
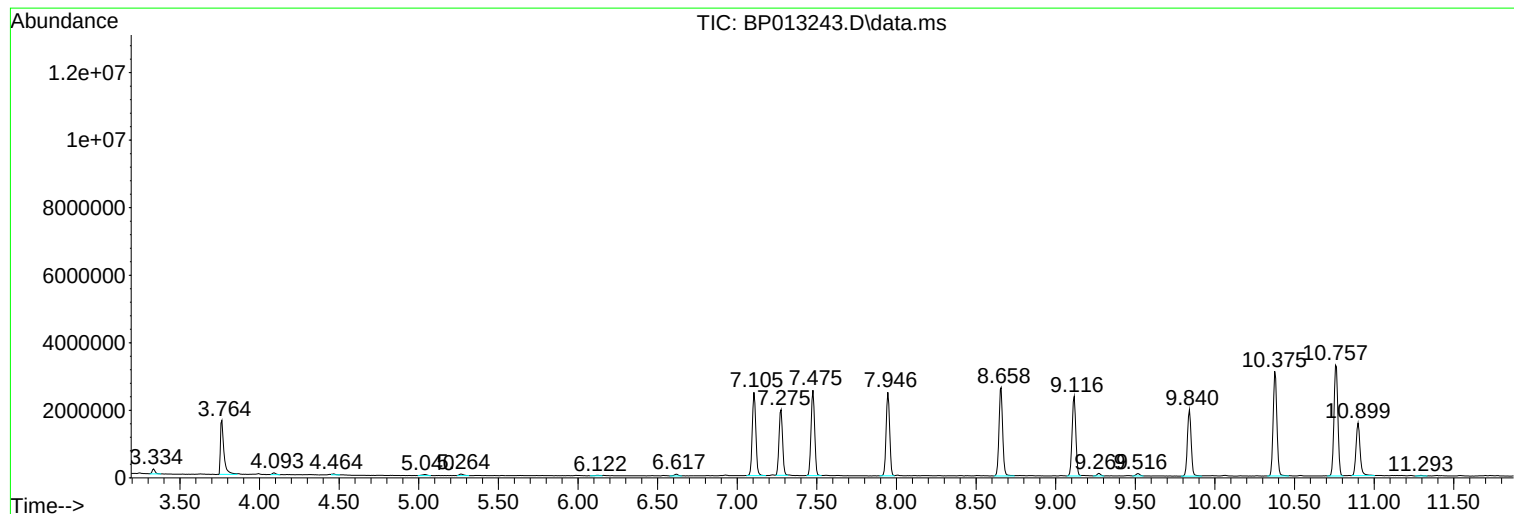
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Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP12222\
Data File : BP013243.D
Acq On : 22 Dec 2022 13:22
Operator : CG/JU
Sample : N6015-02
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DBXT6

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



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122222\
Data File : BP013243.D
Acq On : 22 Dec 2022 13:22
Operator : CG/JU
Sample : N6015-02
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DBXT6

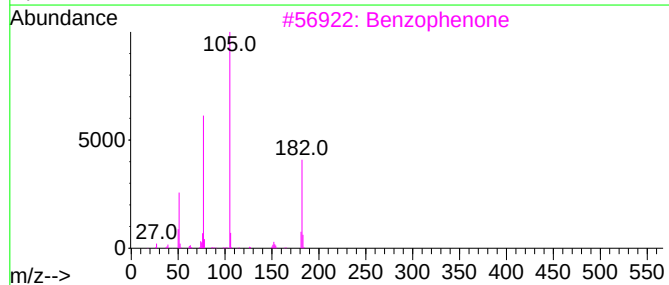
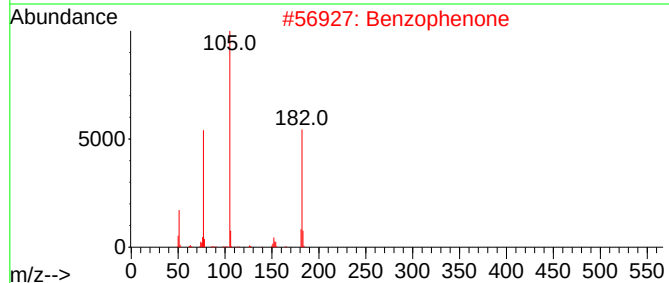
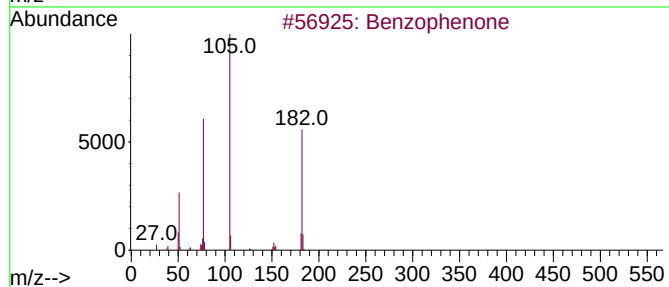
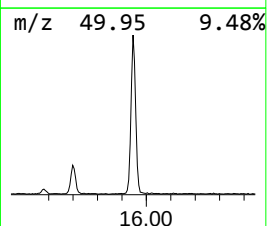
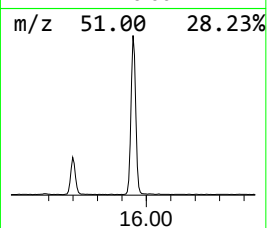
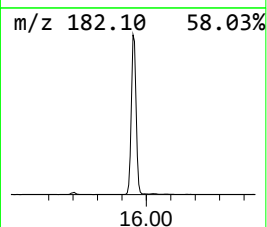
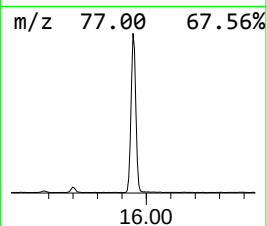
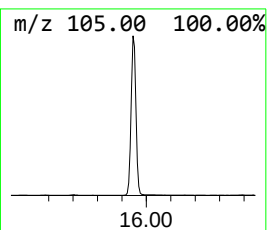
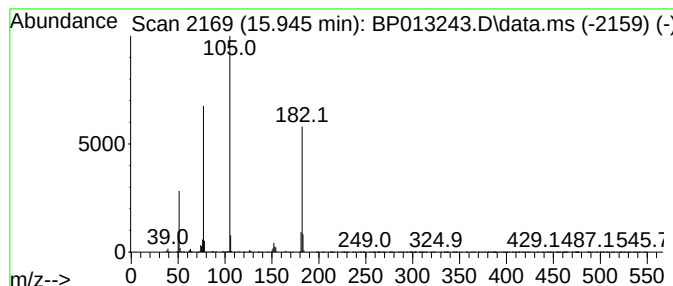
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Benzophenone Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.945	9.55 ng/ul	3557130	Acenaphthene-d10	14.587

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	90
5			Benzenecarbothioic acid, S-methy...	152	C8H8OS	005925-68-8	58



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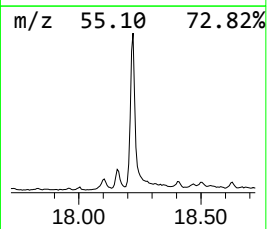
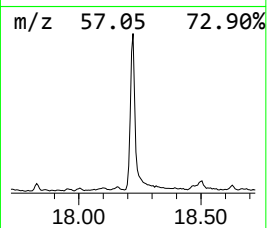
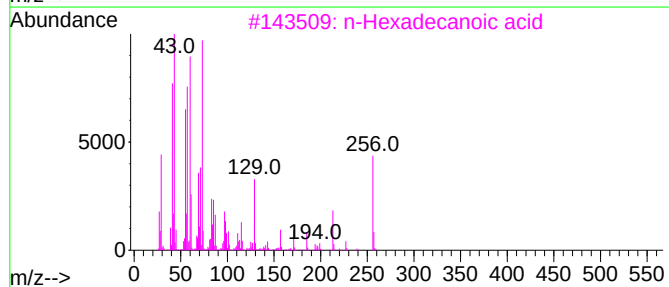
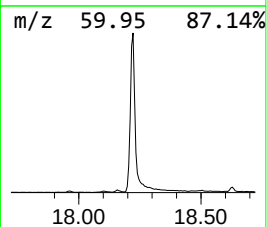
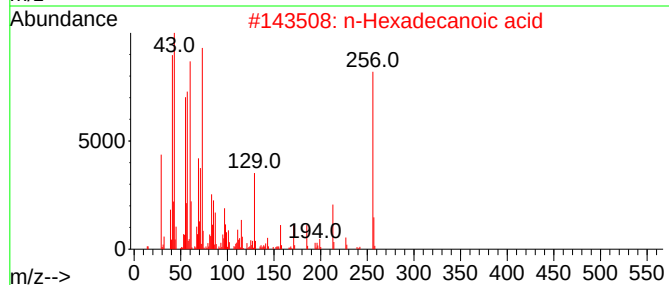
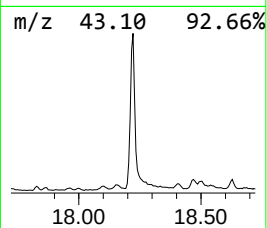
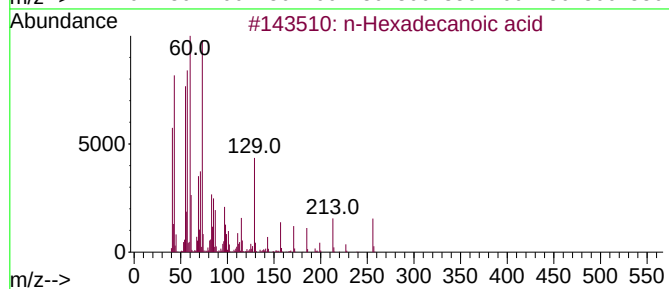
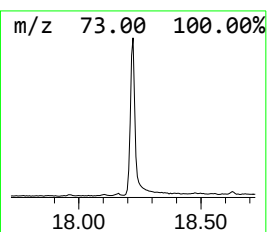
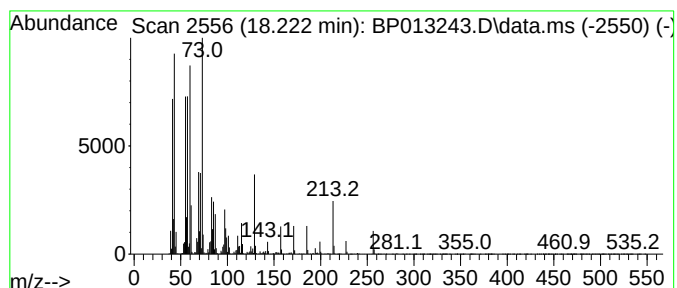
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 n-Hexadecanoic acid Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.222	10.97 ng/ul	4670050	Phenanthrene-d10	17.345

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	98
5			Tridecanoic acid	214	C13H26O2	000638-53-9	95



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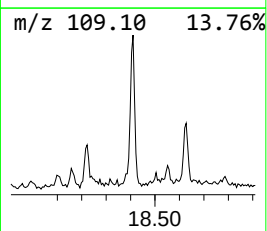
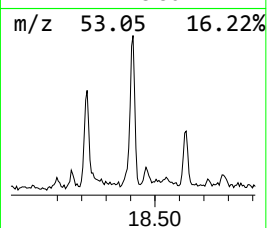
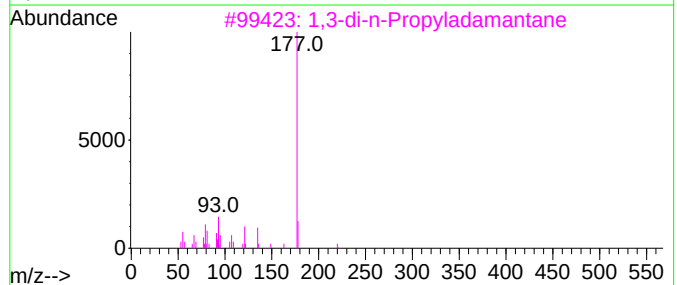
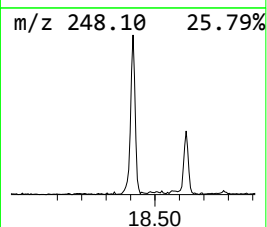
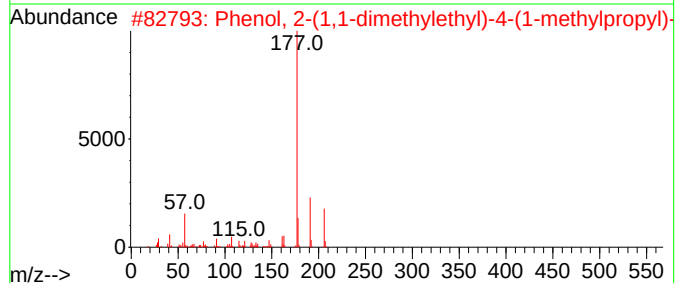
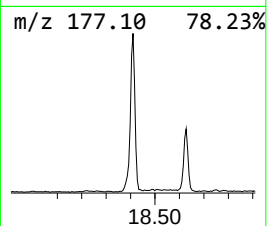
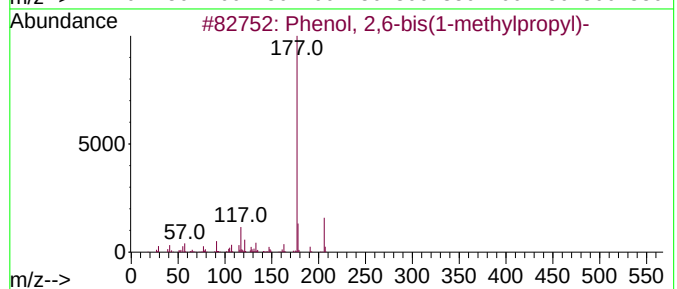
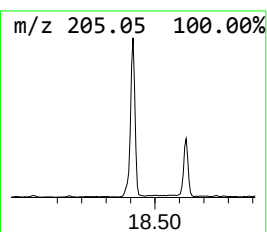
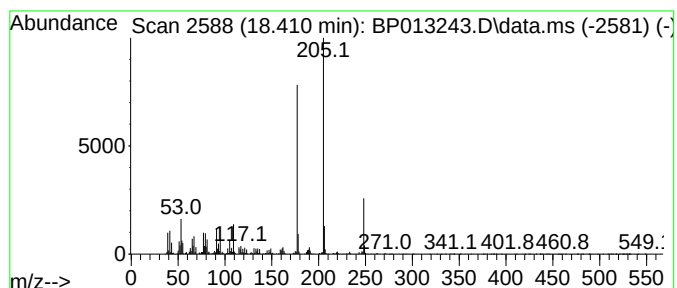
Instrument :
BNA_P
ClientSampleId :
DBXT6

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 unknown-01 Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.	
18.410	2.22 ng/ul	943869	Phenanthrene-d10	17.345	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, 2,6-bis(1-methylpropyl)-	206	C14H22O	005510-99-6	41
2	Phenol, 2-(1,1-dimethylethyl)-4-...	206	C14H22O	052184-13-1	38
3	1,3-di-n-Propyladamantane	220	C16H28	040002-47-9	38
4	Phenol, 2,5-bis(1-methylpropyl)-	206	C14H22O	054932-77-3	38
5	4-Phenylisoquinoline	205	C15H11N	019571-30-3	35



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122222\
Data File : BP013243.D
Acq On : 22 Dec 2022 13:22
Operator : CG/JU
Sample : N6015-02
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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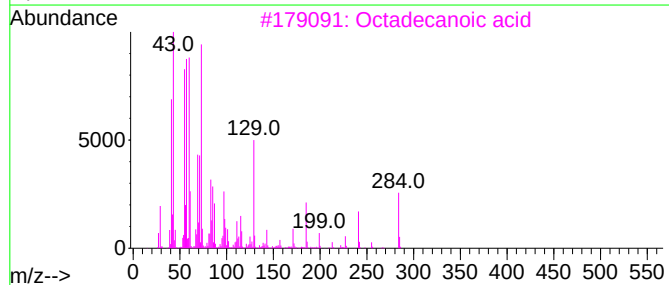
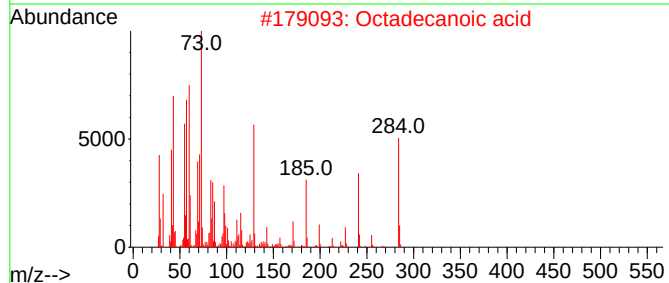
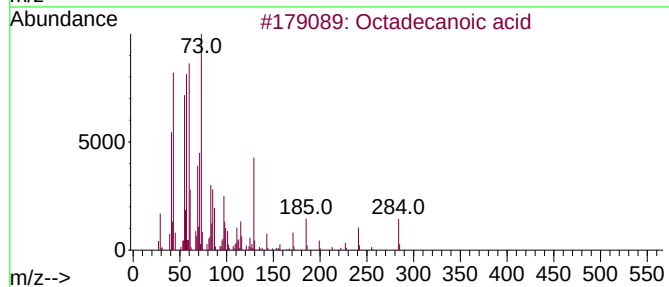
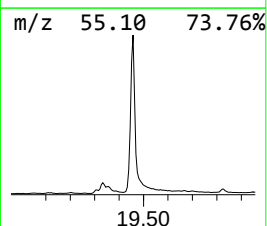
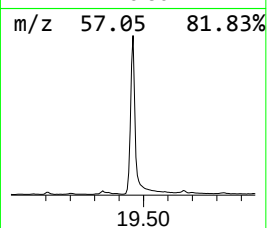
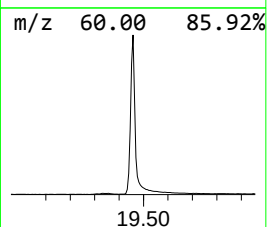
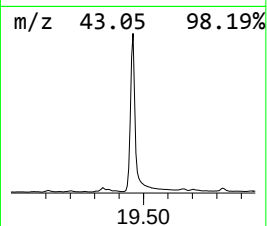
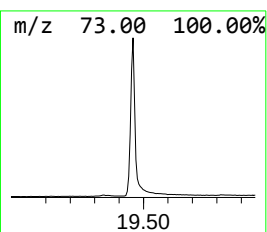
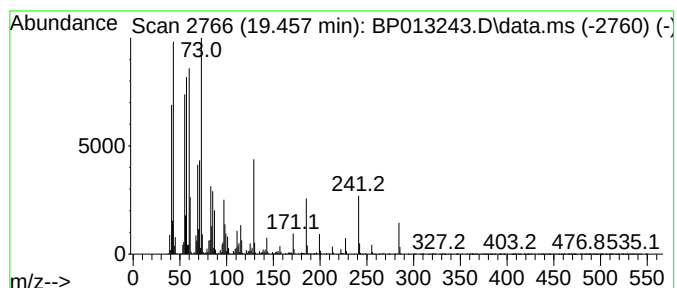
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 Octadecanoic acid Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.457	50.89 ng/ul	17621900	Chrysene-d12	21.433

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octadecanoic acid	284	C18H36O2	000057-11-4	99
2			Octadecanoic acid	284	C18H36O2	000057-11-4	99
3			Octadecanoic acid	284	C18H36O2	000057-11-4	99
4			Octadecanoic acid	284	C18H36O2	000057-11-4	98
5			Octadecanoic acid	284	C18H36O2	000057-11-4	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122222\
Data File : BP013243.D
Acq On : 22 Dec 2022 13:22
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Sample : N6015-02
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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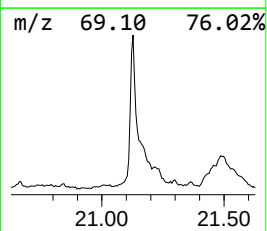
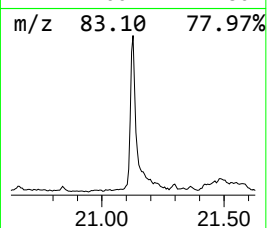
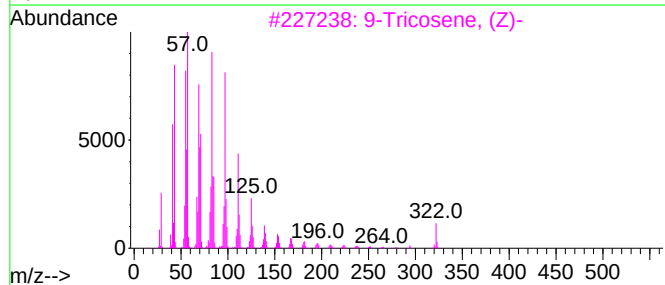
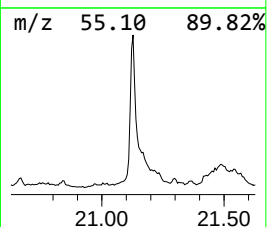
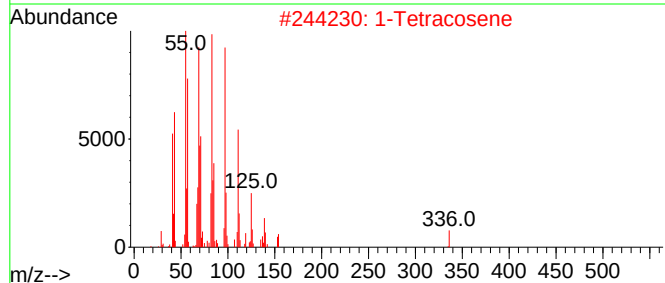
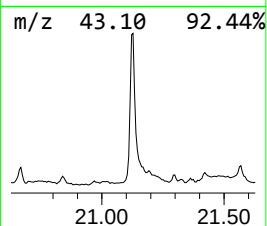
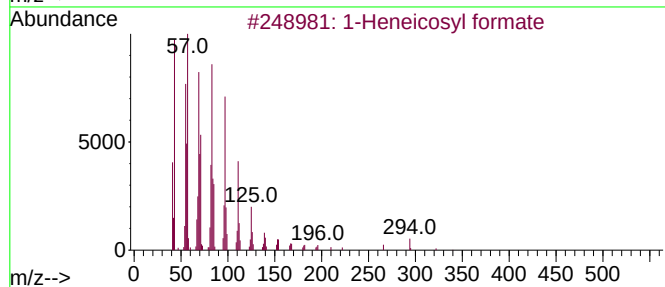
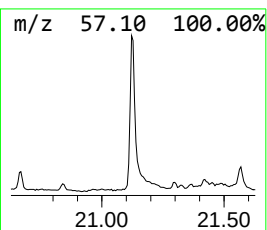
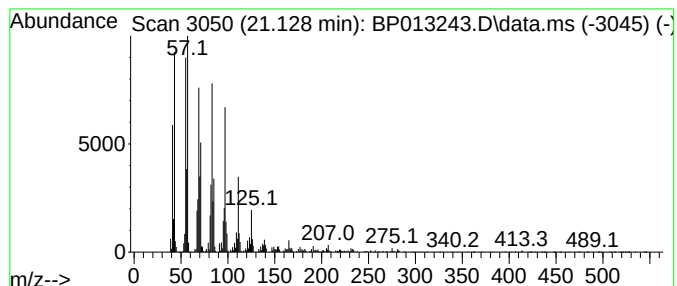
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 1-Heneicosyl formate Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.128	16.97 ng/ul	5876620	Chrysene-d12	21.433

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Heneicosyl formate	340	C22H44O2	077899-03-7	96
2			1-Tetracosene	336	C24H48	010192-32-2	96
3			9-Tricosene, (Z)-	322	C23H46	027519-02-4	95
4			1-Hexacosene	364	C26H52	018835-33-1	94
5			n-Tetracosanol-1	354	C24H50O	000506-51-4	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122222\
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Operator : CG/JU
Sample : N6015-02
Misc :
ALS Vial : 10 Sample Multiplier: 1

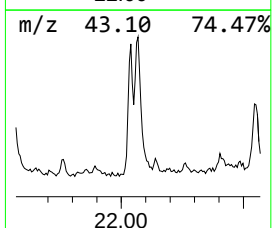
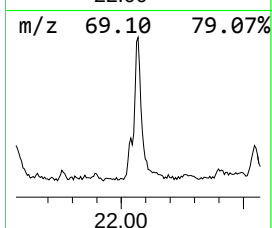
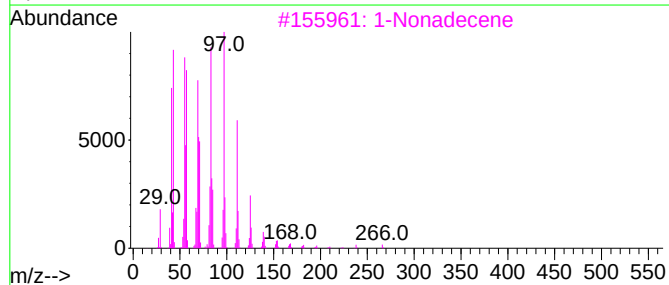
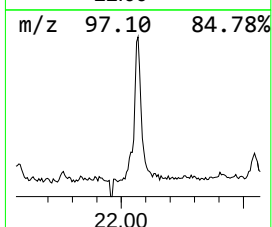
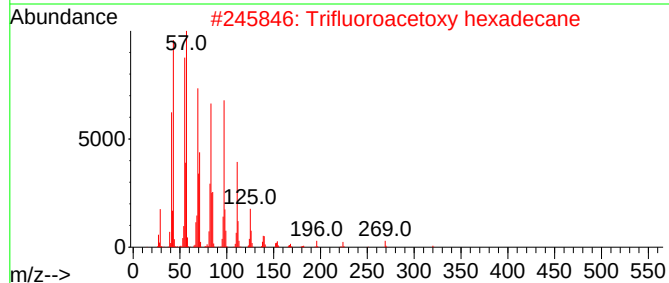
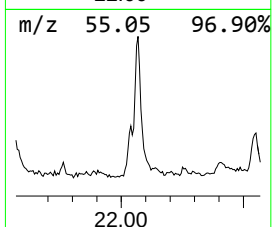
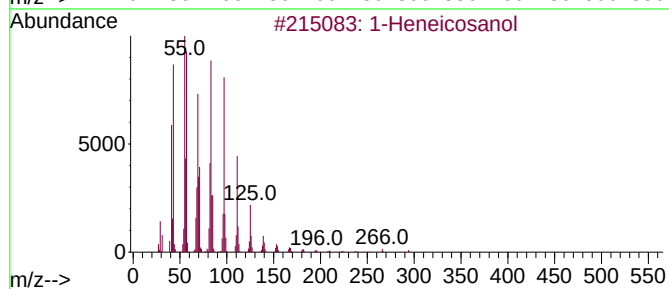
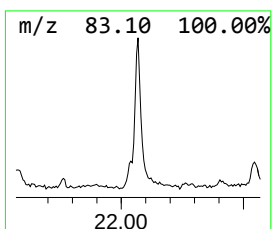
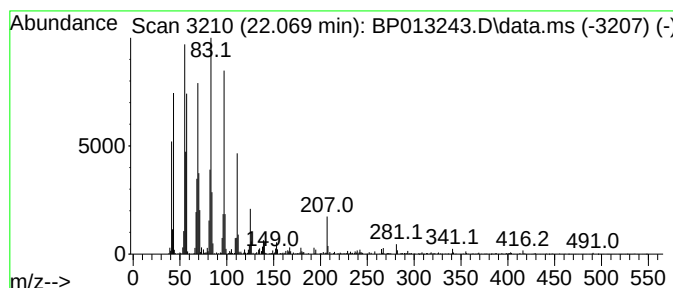
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 1-Heneicosanol Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.	
22.069	3.45 ng/ul	1194550	Chrysene-d12	21.433	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Heneicosanol	312	C21H44O	015594-90-8	94
2	Trifluoroacetoxy hexadecane	338	C18H33F3O2	006222-03-3	91
3	1-Nonadecene	266	C19H38	018435-45-5	90
4	Behenic alcohol	326	C22H46O	000661-19-8	90
5	1-Docosene	308	C22H44	001599-67-3	89



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122222\
Data File : BP013243.D
Acq On : 22 Dec 2022 13:22
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Sample : N6015-02
Misc :
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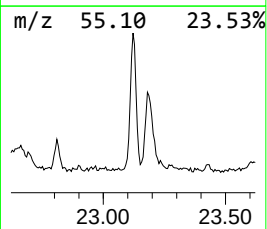
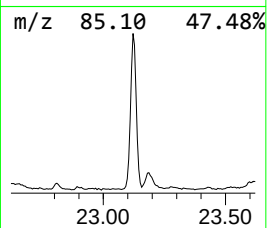
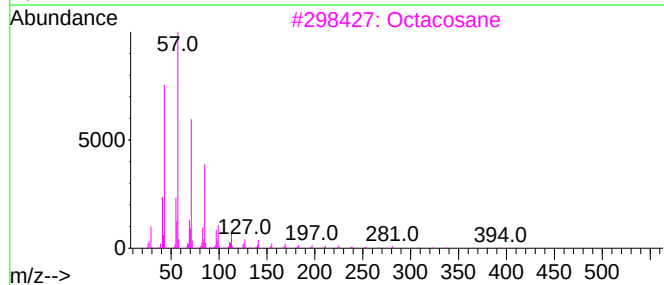
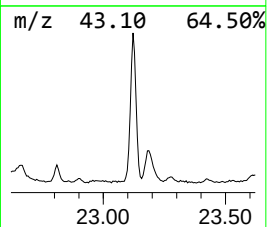
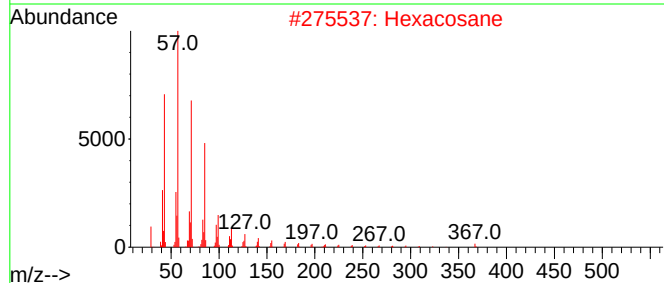
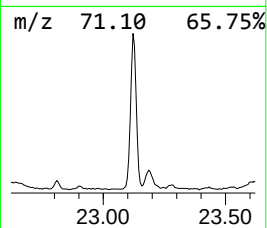
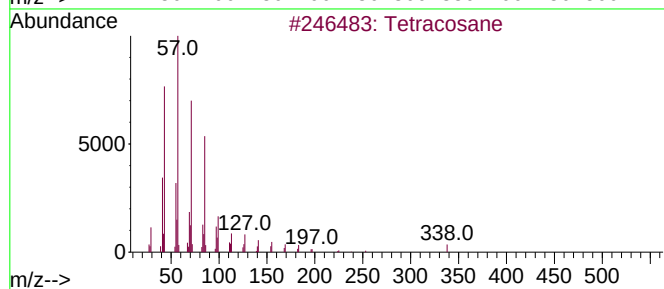
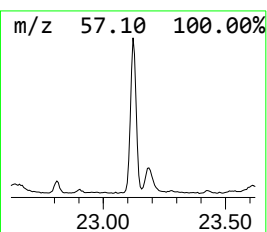
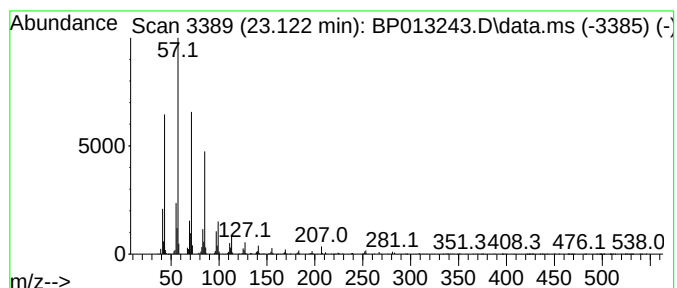
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 (DEL) Alkane: Straight-Chai... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.122	5.36 ng/ul	1793620	Perylene-d12	23.910

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tetracosane	338	C24H50	000646-31-1	93
2			Hexacosane	366	C26H54	000630-01-3	90
3			Octacosane	394	C28H58	000630-02-4	90
4			Heneicosane	296	C21H44	000629-94-7	87
5			Tetracosane	338	C24H50	000646-31-1	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122222\
Data File : BP013243.D
Acq On : 22 Dec 2022 13:22
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Sample : N6015-02
Misc :
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Instrument :
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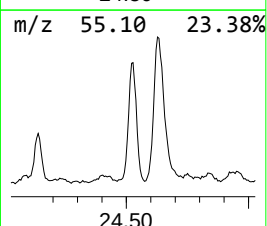
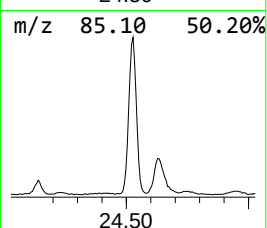
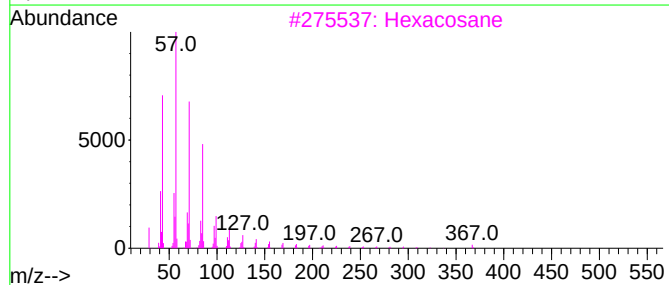
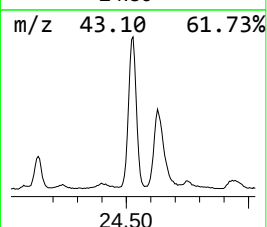
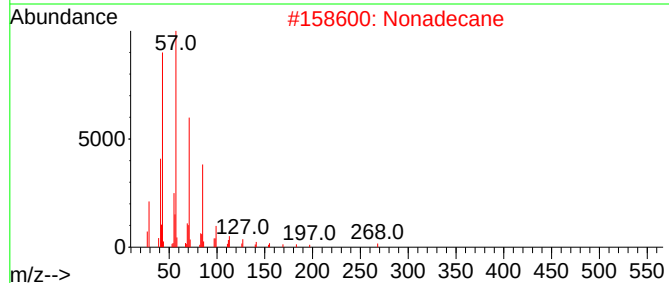
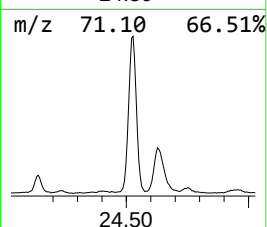
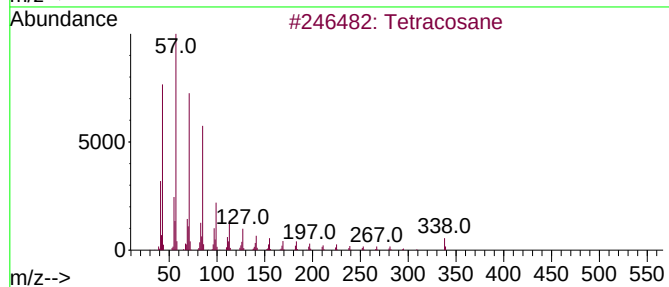
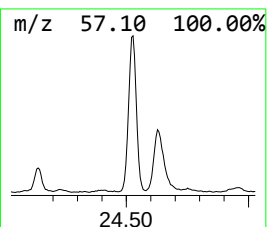
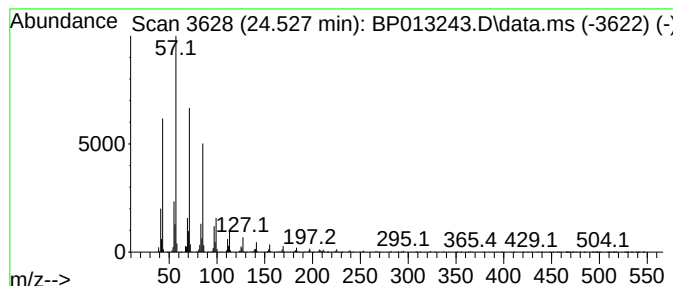
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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.527	12.51 ng/ul	4188070	Perylene-d12	23.910

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Tetracosane	338	C24H50	000646-31-1	96
2		Nonadecane	268	C19H40	000629-92-5	95
3		Hexacosane	366	C26H54	000630-01-3	94
4		2-Methyltetracosane	352	C25H52	001560-78-7	94
5		Octadecane, 2-methyl-	268	C19H40	001560-88-9	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122222\
Data File : BP013243.D
Acq On : 22 Dec 2022 13:22
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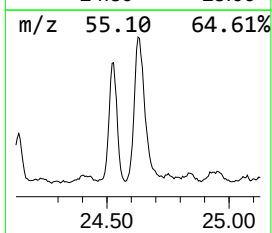
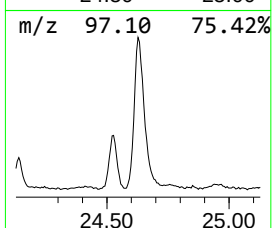
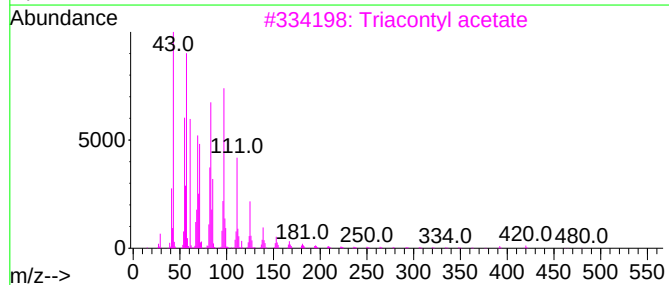
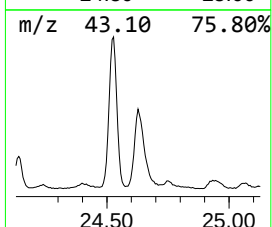
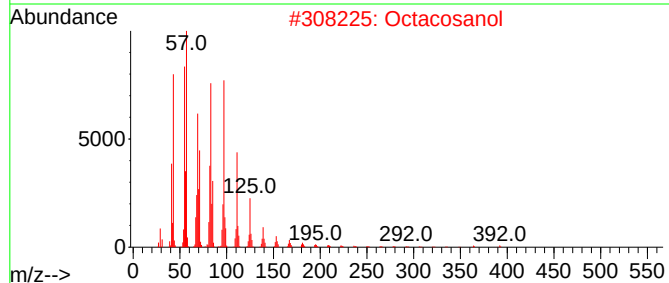
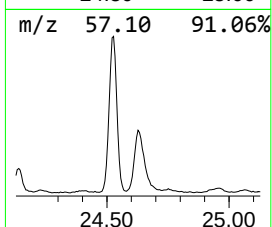
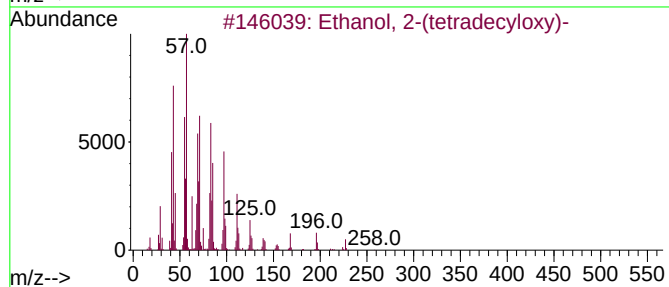
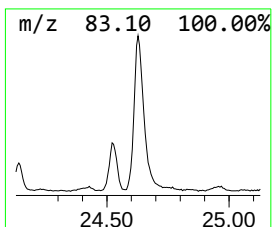
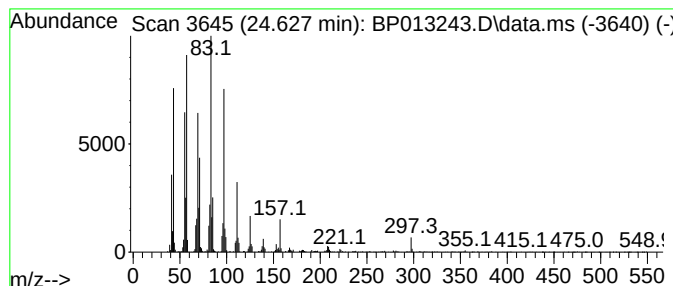
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Quant Title : SVOA CALIBRATION

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

Peak Number 11 Ethanol, 2-(tetradecyloxy)- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.627	12.15 ng/ul	4065950	Perylene-d12	23.910

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Ethanol, 2-(tetradecyloxy)-	258	C16H34O2	002136-70-1	83
2			Octacosanol	410	C28H58O	000557-61-9	60
3			Triacontyl acetate	480	C32H64O2	041755-58-2	60
4			Heptafluorobutyric acid, hexadec...	438	C20H33F7O2	006385-15-5	58
5			Triacontan-15-ol	438	C30H62O	031849-26-0	58



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122222\
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Sample : N6015-02
Misc :
ALS Vial : 10 Sample Multiplier: 1

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ClientSampleId :
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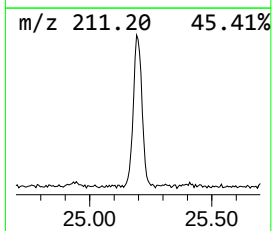
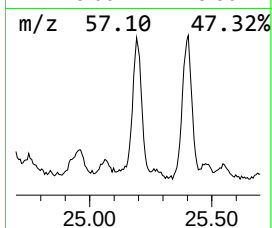
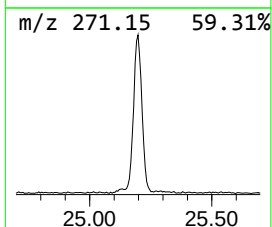
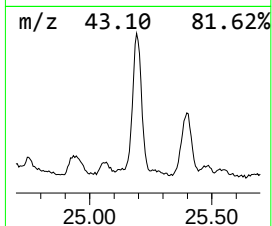
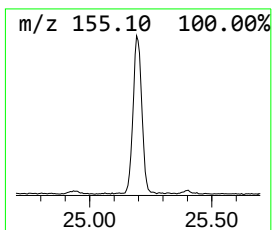
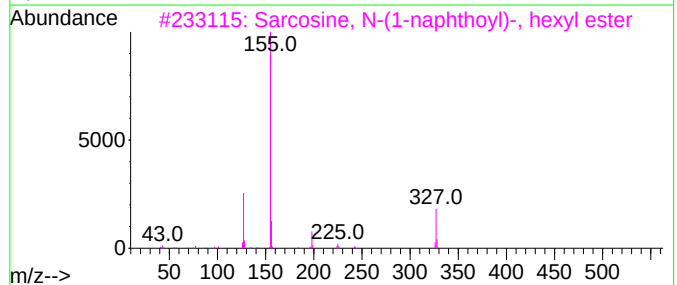
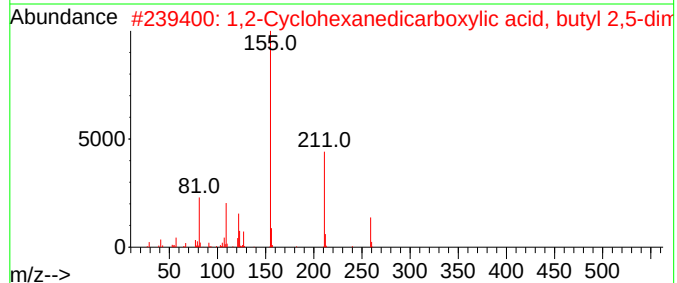
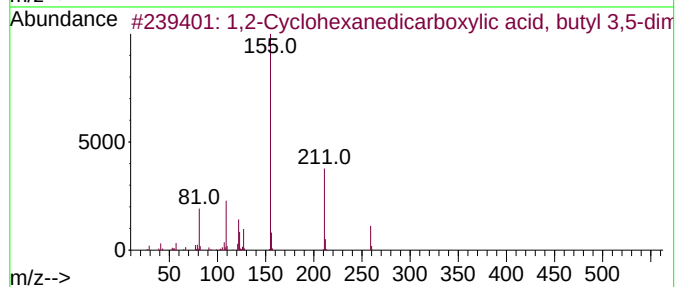
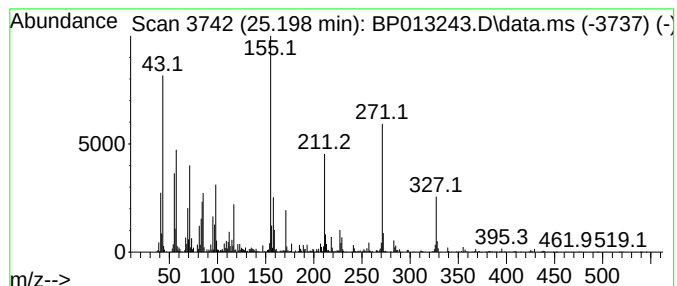
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Quant Title : SVOA CALIBRATION

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

Peak Number 12 unknown-02 Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
25.198	7.26 ng/ul	2428440	Perylene-d12	23.910

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1,2-Cyclohexanedicarboxylic acid...	332	C20H28O4	1000339-61-4	35
2			1,2-Cyclohexanedicarboxylic acid...	332	C20H28O4	1000339-94-0	35
3			Sarcosine, N-(1-naphthoyl)-, hex...	327	C20H25NO3	1000321-40-4	30
4			4-(Decanoyloxy)benzoic acid	292	C17H24O4	086960-46-5	27
5			Sarcosine, N-(1-naphthoyl)-, iso...	327	C20H25NO3	1010321-40-3	25



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122222\
Data File : BP013243.D
Acq On : 22 Dec 2022 13:22
Operator : CG/JU
Sample : N6015-02
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DBXT6

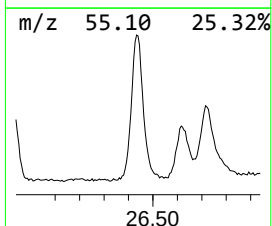
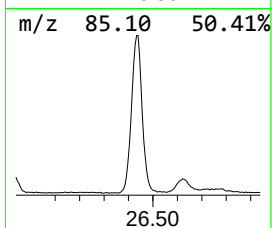
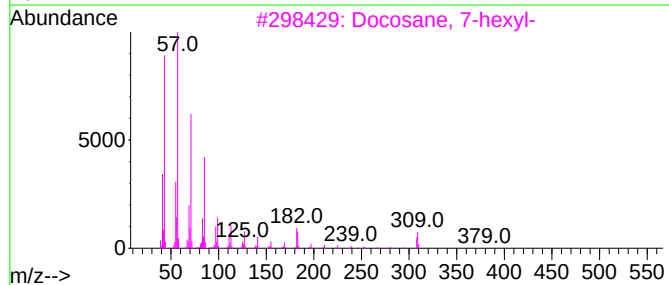
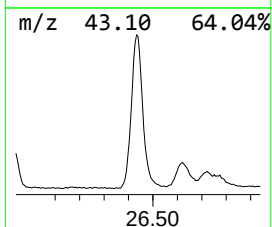
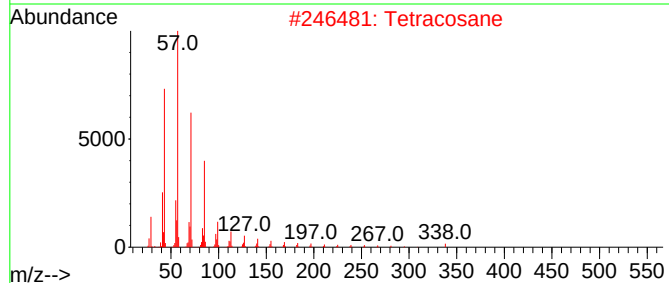
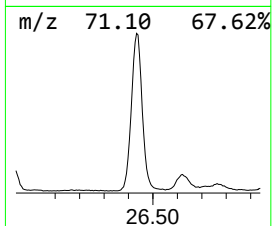
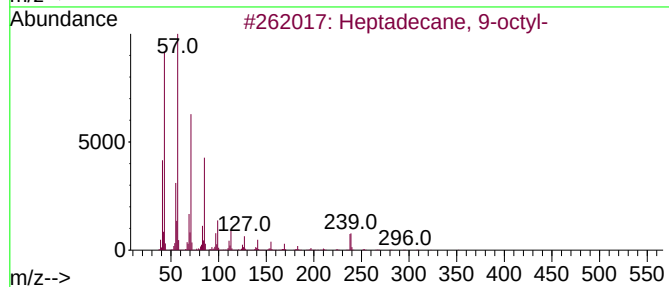
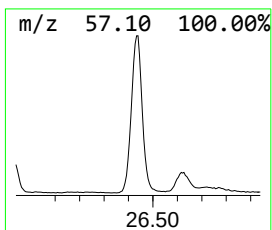
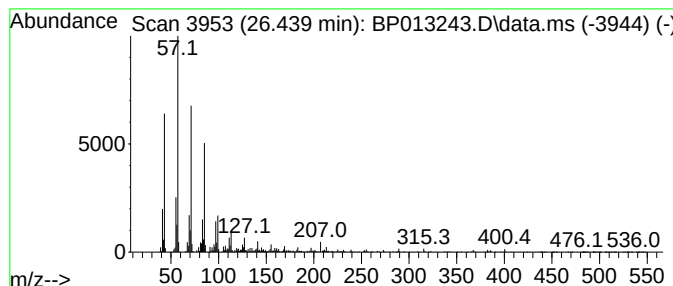
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Quant Title : SVOA CALIBRATION

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

Peak Number 13 (DEL) Alkane: Straight-Chai... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
26.439	26.99 ng/ul	9034820	Perylene-d12	23.910

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Heptadecane, 9-octyl-	352	C25H52	007225-64-1	94
2		Tetracosane	338	C24H50	000646-31-1	93
3		Docosane, 7-hexyl-	394	C28H58	055373-86-9	93
4		Octadecane, 5,14-dibutyl-	366	C26H54	055282-13-8	93
5		2-Methylpentacosane	366	C26H54	000629-87-8	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122222\
Data File : BP013243.D
Acq On : 22 Dec 2022 13:22
Operator : CG/JU
Sample : N6015-02
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DBXT6

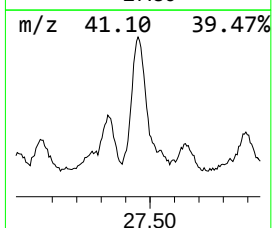
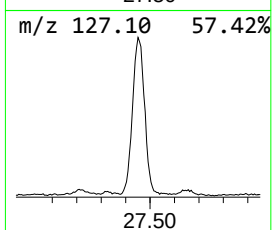
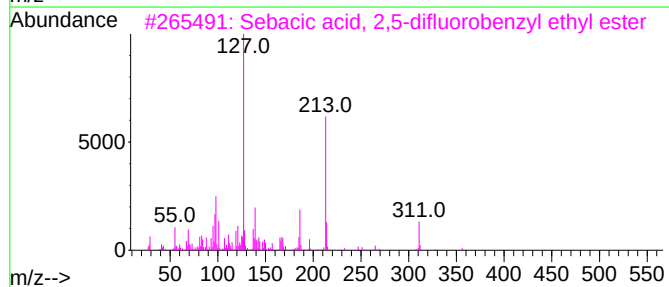
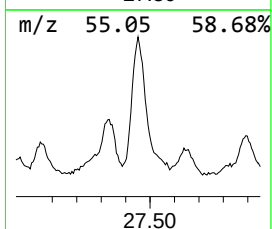
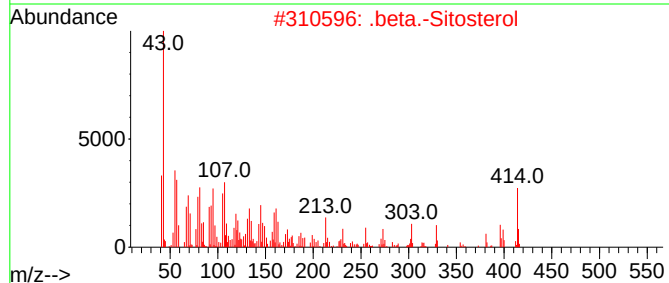
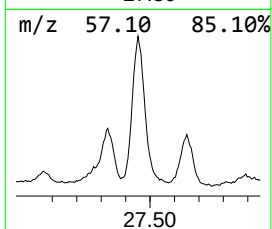
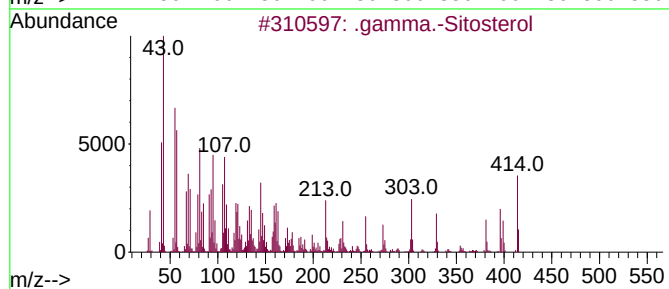
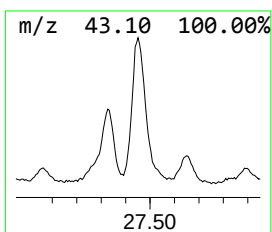
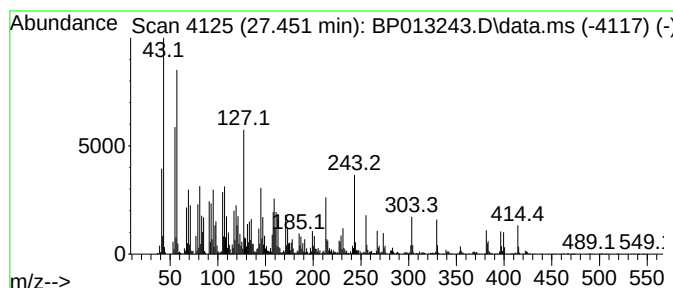
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Quant Title : SVOA CALIBRATION

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

Peak Number 14 .gamma.-Sitosterol Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
27.451	20.04 ng/ul	6707260	Perylene-d12	23.910

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	.	gamma.-Sitosterol	414	C29H50O	000083-47-6	97	
2	.	beta.-Sitosterol	414	C29H50O	000083-46-5	58	
3		Sebacic acid, 2,5-difluorobenzyl...	356	C19H26F2O4	1000380-76-0	38	
4		(3S,8S,9S,10R,13R,14S,17R)-17-((...	414	C29H50O	029365-29-5	15	
5		5-(4-Chloroanilino)-5-oxopentano...	241	C11H12ClNO3	040828-92-0	11	



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122222\
Data File : BP013243.D
Acq On : 22 Dec 2022 13:22
Operator : CG/JU
Sample : N6015-02
Misc :
ALS Vial : 10 Sample Multiplier: 1

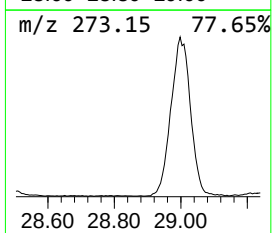
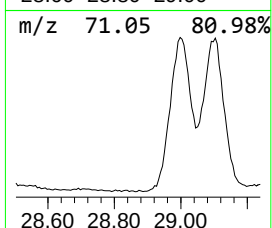
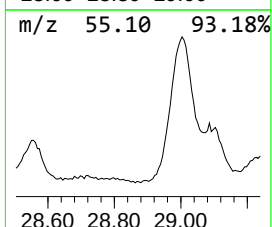
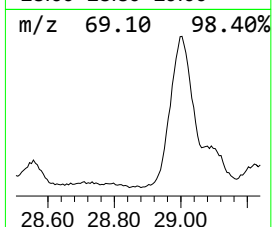
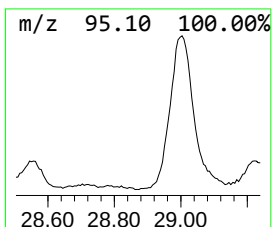
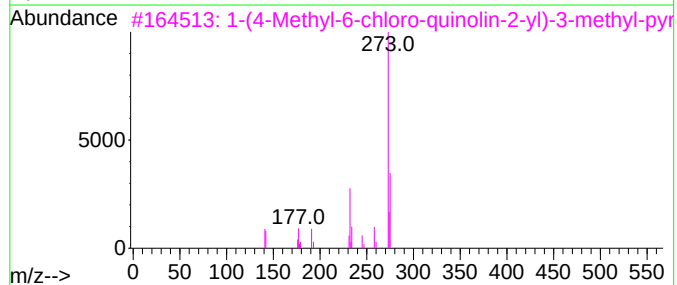
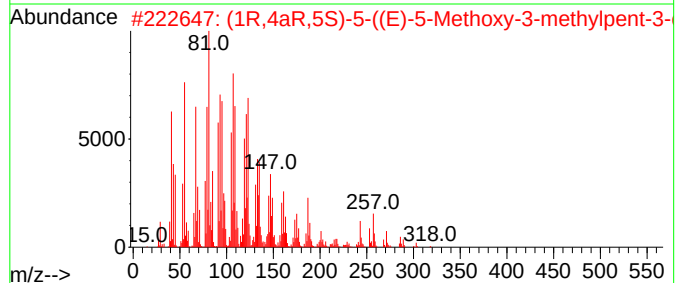
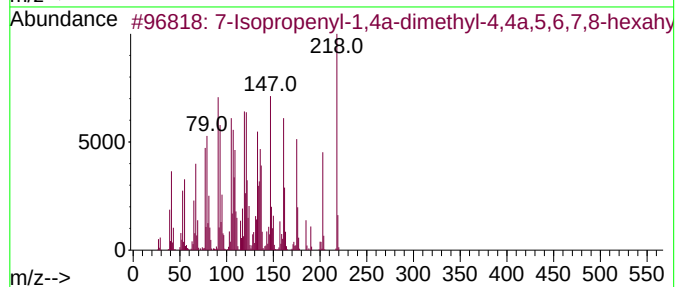
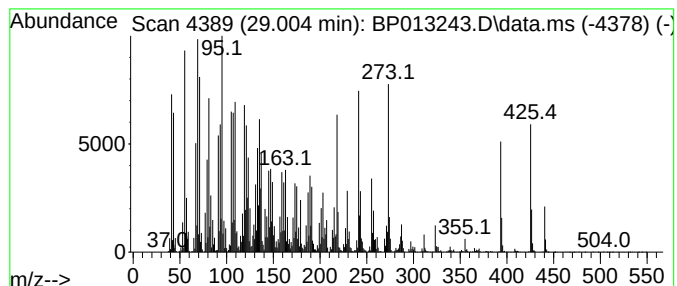
Instrument :
BNA_P
ClientSampleId :
DBXT6

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 15 unknown-03 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD		R.T.	
29.004	62.26 ng/ul	20836400	Perylene-d12		23.910	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	7-Isopropenyl-1,4a-dimethyl-4,4a...		218	C15H22O	000473-08-5	40
2	(1R,4aR,5S)-5-((E)-5-Methoxy-3-m...		318	C21H34O2	1000495-78-7	25
3	1-(4-Methyl-6-chloro-quinolin-2-...		273	C14H12ClN3O	111784-26-0	25
4	D:C-Friedo-B':A'-neogammacer-9(1...		440	C31H52O	004555-56-0	22
5	9-Thiabicyclo[3.3.1]nonane, 2,6-...		274	C14H18N4S	1000141-52-1	22



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122222\
 Data File : BP013243.D
 Acq On : 22 Dec 2022 13:22
 Operator : CG/JU
 Sample : N6015-02
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 DBXT6

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
Benzophenone	15.945	9.6	ng/u1	3557130	3	14.587	7447470	20.0
n-Hexadecanoic ...	18.222	11.0	ng/u1	4670050	4	17.345	8512620	20.0
unknown-01	18.410	2.2	ng/u1	943869	4	17.345	8512620	20.0
Octadecanoic acid	19.457	50.9	ng/u1	17621900	5	21.433	6925150	20.0
4-(3-Fluoro-8-n...	19.828	12.5	ng/u1	4341460	5	21.433	6925150	20.0
(1H)Pyrrole, 3,...	20.169	3.0	ng/u1	1031340	5	21.433	6925150	20.0
1-Heneicosyl fo...	21.128	17.0	ng/u1	5876620	5	21.433	6925150	20.0
1-Heneicosanol	22.069	3.5	ng/u1	1194550	5	21.433	6925150	20.0
(DEL) Alkane: S...	23.122	5.4	ng/u1	1793620	6	23.910	6693720	20.0
(DEL) Alkane: S...	24.527	12.5	ng/u1	4188070	6	23.910	6693720	20.0
Ethanol, 2-(tet...	24.627	12.2	ng/u1	4065950	6	23.910	6693720	20.0
unknown-02	25.198	7.3	ng/u1	2428440	6	23.910	6693720	20.0
(DEL) Alkane: S...	26.439	27.0	ng/u1	9034820	6	23.910	6693720	20.0
.gamma.-Sitosterol	27.451	20.0	ng/u1	6707260	6	23.910	6693720	20.0
unknown-03	29.004	62.3	ng/u1	20836400	6	23.910	6693720	20.0