

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP12222\
 Data File : BP013246.D
 Acq On : 22 Dec 2022 15:09
 Operator : CG/JU
 Sample : N6015-05
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 DBXU1

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: BP013246.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	29	rBV	137870	168949	1.47%	0.102%
2	3.764	95	98	113	rBV	1716810	2566208	22.30%	1.553%
3	3.870	113	116	122	rVB2	36609	52484	0.46%	0.032%
4	3.993	135	137	147	rVB4	25044	37622	0.33%	0.023%
5	4.087	150	153	158	rVB	48458	65458	0.57%	0.040%
6	5.034	309	314	323	rVB2	44797	81315	0.71%	0.049%
7	5.270	349	354	361	rVB	26937	46659	0.41%	0.028%
8	5.699	422	427	431	rBV6	15104	23433	0.20%	0.014%
9	6.928	627	636	642	rBV3	21780	46584	0.40%	0.028%
10	7.105	660	666	679	rVB	2248837	3539526	30.76%	2.142%
11	7.275	688	695	705	rBV	1778878	2788612	24.24%	1.687%
12	7.475	722	729	740	rBV	2308900	3556063	30.91%	2.152%
13	7.946	802	809	816	rBV	2241415	3451069	29.99%	2.088%
14	8.005	816	819	823	rVB3	17846	25214	0.22%	0.015%
15	8.658	922	930	942	rBV	2528902	4096266	35.60%	2.479%
16	9.116	1000	1008	1016	rBV	2112882	3467304	30.13%	2.098%
17	9.269	1030	1034	1043	rVB8	28847	56864	0.49%	0.034%
18	9.516	1070	1076	1081	rBV	61767	96448	0.84%	0.058%
19	9.705	1105	1108	1114	rVB	23888	33732	0.29%	0.020%
20	9.840	1122	1131	1145	rBV	1826758	3004092	26.11%	1.818%
21	10.058	1165	1168	1175	rVB9	12557	24663	0.21%	0.015%
22	10.322	1206	1213	1216	rVV2	54829	87183	0.76%	0.053%
23	10.381	1216	1223	1235	rVB	2729510	4611121	40.08%	2.790%
24	10.763	1278	1288	1294	rBV	2931898	4970604	43.20%	3.008%
25	10.899	1300	1311	1331	rVV	1574299	2841523	24.70%	1.719%
26	11.287	1373	1377	1382	rVB8	13315	22706	0.20%	0.014%
27	11.769	1457	1459	1466	rVB8	15410	23304	0.20%	0.014%
28	12.081	1506	1512	1515	rBV3	16361	24886	0.22%	0.015%
29	12.169	1522	1527	1532	rBV2	38010	58368	0.51%	0.035%
30	12.252	1535	1541	1545	rBV5	12787	23048	0.20%	0.014%
31	12.340	1552	1556	1562	rVV2	46981	79846	0.69%	0.048%
32	12.416	1566	1569	1574	rVB3	16847	24757	0.22%	0.015%
33	12.557	1583	1593	1600	rBV3	8054	27000	0.23%	0.016%
34	13.116	1684	1688	1694	rVB5	23385	36547	0.32%	0.022%
35	13.322	1718	1723	1726	rBV6	16771	25173	0.22%	0.015%
36	13.404	1728	1737	1744	rVB	79406	132864	1.15%	0.080%

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Integration Parameters: LSCINT.P

Integrator: RTE

Smoothing : OFF

Filtering: 5

Sampling : 1

Min Area: 0.1 % of largest Peak

Start Thrs: 0.2

Max Peaks: 100

Stop Thrs : 0

Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >

Peak separation: 5

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

37	13.516	1751	1756	1759	rBV5	17306	27999	0.24%	0.017%
38	13.769	1795	1799	1808	rVB10	14644	29138	0.25%	0.018%
39	13.910	1818	1823	1829	rBV7	14387	32187	0.28%	0.019%
40	13.993	1829	1837	1845	rBV	4225701	5965388	51.85%	3.610%
41	14.057	1846	1848	1852	rVB3	26341	33392	0.29%	0.020%
42	14.104	1852	1856	1861	rVB7	19195	26007	0.23%	0.016%
43	14.181	1865	1869	1872	rVB2	19717	24712	0.21%	0.015%
44	14.281	1879	1886	1898	rBV	4937571	7347200	63.85%	4.446%
45	14.475	1916	1919	1924	rVB2	50899	63975	0.56%	0.039%
46	14.587	1930	1938	1948	rBV	4087836	6009750	52.23%	3.637%
47	14.787	1965	1972	1994	rBV	1334182	2216744	19.27%	1.341%
48	14.946	1994	1999	2002	rVV5	37256	63629	0.55%	0.039%
49	14.993	2002	2007	2013	rVV4	55868	99905	0.87%	0.060%
50	15.369	2067	2071	2074	rBV5	17862	27845	0.24%	0.017%
51	15.422	2078	2080	2086	rVB2	23708	30429	0.26%	0.018%
52	15.581	2100	2107	2121	rBV	6022651	8713673	75.73%	5.273%
53	15.698	2122	2127	2136	rBV	1548459	2116869	18.40%	1.281%
54	15.822	2145	2148	2155	rVB4	32104	53933	0.47%	0.033%
55	15.945	2159	2169	2177	rBV	2343504	3346448	29.08%	2.025%
56	16.145	2197	2203	2210	rVB	19587	40770	0.35%	0.025%
57	16.216	2210	2215	2221	rBV4	38045	57102	0.50%	0.035%
58	16.310	2224	2231	2236	rVV6	31123	73695	0.64%	0.045%
59	16.434	2249	2252	2253	rBV3	22085	22598	0.20%	0.014%
60	16.516	2262	2266	2275	rVB2	82987	119359	1.04%	0.072%
61	16.728	2298	2302	2310	rVB5	63343	100519	0.87%	0.061%
62	16.945	2336	2339	2344	rBV	59113	86995	0.76%	0.053%
63	17.087	2361	2363	2367	rVB4	21707	23908	0.21%	0.014%
64	17.228	2379	2387	2394	rBV3	109361	195270	1.70%	0.118%
65	17.292	2394	2398	2401	rVV2	60895	70649	0.61%	0.043%
66	17.345	2401	2407	2417	rVV	4136767	6127001	53.25%	3.708%
67	17.445	2417	2424	2431	rVB	5477307	7727713	67.16%	4.676%
68	17.557	2440	2443	2450	rVB5	43471	53334	0.46%	0.032%
69	17.622	2451	2454	2463	rVB8	28479	52911	0.46%	0.032%
70	17.698	2463	2467	2480	rBV5	86029	159248	1.38%	0.096%
71	17.863	2492	2495	2500	rVB4	24137	41384	0.36%	0.025%
72	18.004	2516	2519	2523	rVB4	39186	41796	0.36%	0.025%
73	18.104	2531	2536	2541	rBV5	87567	146558	1.27%	0.089%
74	18.157	2541	2545	2550	rBV2	189775	268166	2.33%	0.162%
75	18.216	2550	2555	2574	rVV	1709712	2600504	22.60%	1.574%
76	18.410	2584	2588	2593	rVB	136502	186535	1.62%	0.113%

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Integration Parameters: LSCINT.P

Integrator: RTE
 Smoothing : OFF Filtering: 5
 Sampling : 1 Min Area: 0.1 % of largest Peak
 Start Thrs: 0.2 Max Peaks: 100
 Stop Thrs : 0 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
 Peak separation: 5

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77	18.463	2594	2597	2600	rBV	105272	121199	1.05%	0.073%
78	18.504	2600	2604	2608	rVV4	67702	94898	0.82%	0.057%
79	18.628	2622	2625	2630	rBV5	83730	120676	1.05%	0.073%
80	18.781	2647	2651	2659	rVB	280115	374375	3.25%	0.227%
81	19.122	2704	2709	2713	rBV2	225069	300179	2.61%	0.182%
82	19.204	2719	2723	2728	rBV2	106993	177624	1.54%	0.107%
83	19.251	2728	2731	2735	rVB3	116024	144532	1.26%	0.087%
84	19.334	2737	2745	2746	rBV3	270278	547616	4.76%	0.331%
85	19.445	2760	2764	2772	rBV	975249	1456536	12.66%	0.881%
86	19.675	2798	2803	2811	rBV	6409084	8499506	73.87%	5.143%
87	19.828	2824	2829	2841	rBV	4441244	5135338	44.63%	3.107%
88	19.951	2847	2850	2855	rBV4	147586	257236	2.24%	0.156%
89	20.004	2857	2859	2864	rVB2	112845	137254	1.19%	0.083%
90	20.169	2883	2887	2894	rVB3	477592	724355	6.30%	0.438%
91	21.128	3045	3050	3053	rBV	1783106	2822577	24.53%	1.708%
92	21.363	3087	3090	3094	rBV3	263065	354314	3.08%	0.214%
93	21.433	3097	3102	3106	rBV	4502981	6162005	53.55%	3.729%
94	23.122	3385	3389	3395	rBV2	804164	1370621	11.91%	0.829%
95	23.751	3489	3496	3510	rVB2	4898118	10201993	88.67%	6.173%
96	23.910	3517	3523	3533	rVB2	3117311	6566758	57.07%	3.974%
97	24.527	3622	3628	3638	rVB	1487899	3247477	28.22%	1.965%
98	24.639	3642	3647	3662	rVB	988222	2614158	22.72%	1.582%
99	26.439	3945	3953	3974	rVB	2190195	7677355	66.72%	4.646%
100	29.004	4379	4389	4402	rBV7	2856903	11506097	100.00%	6.963%

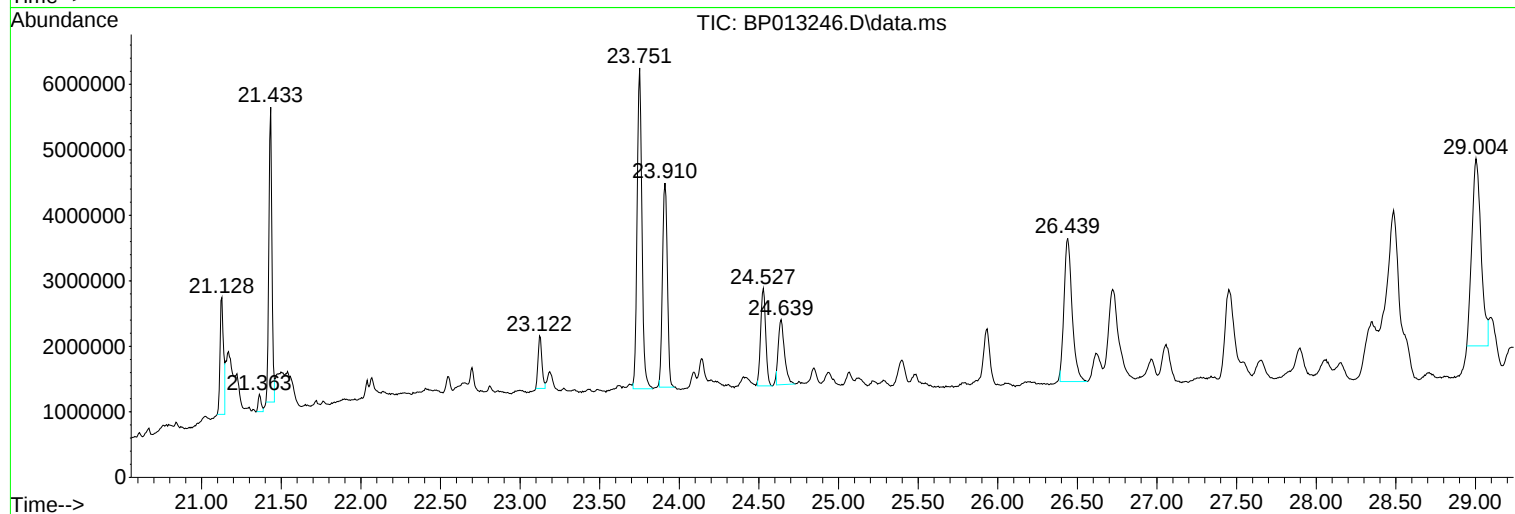
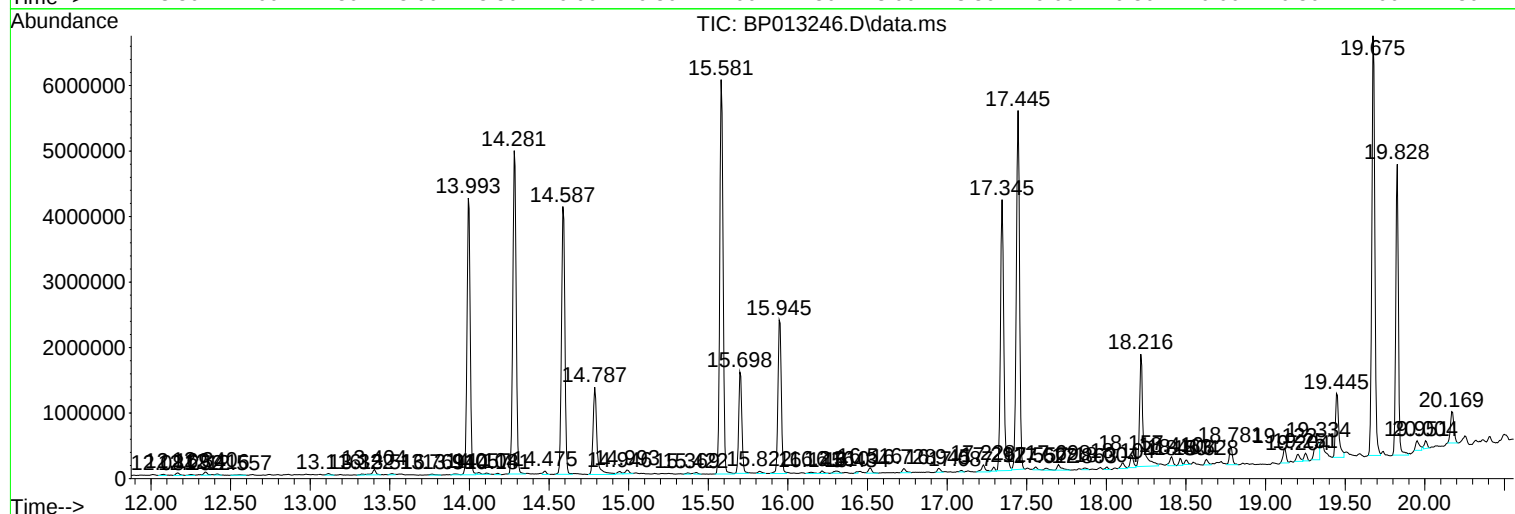
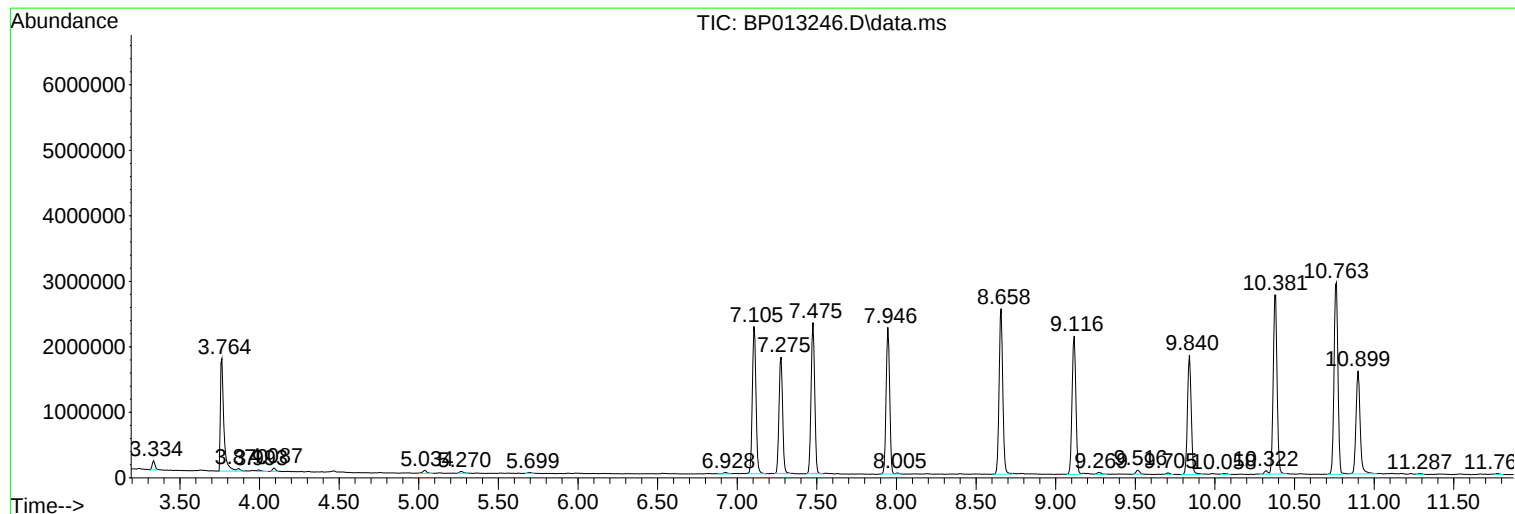
Sum of corrected areas: 165257412

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Instrument :
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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



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122222\
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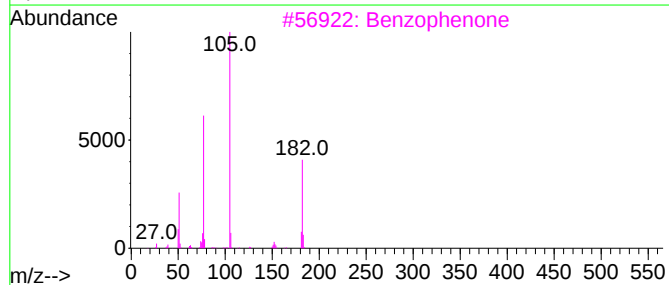
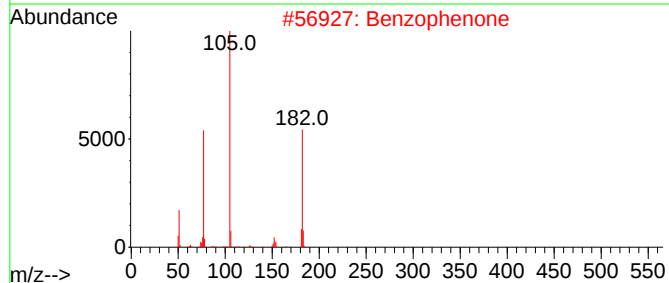
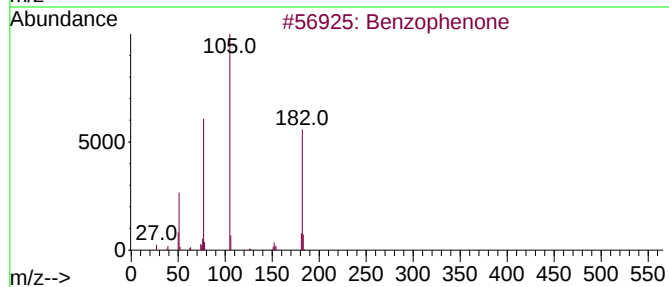
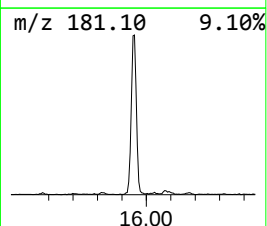
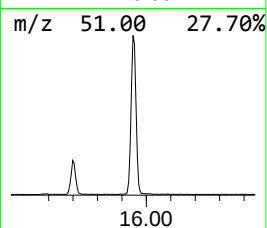
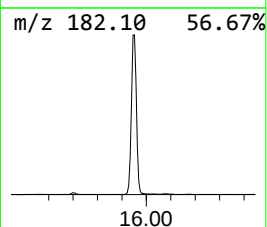
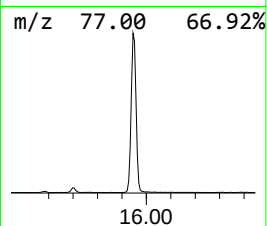
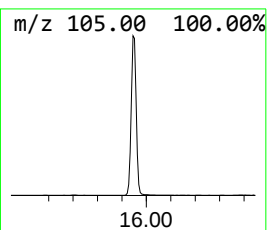
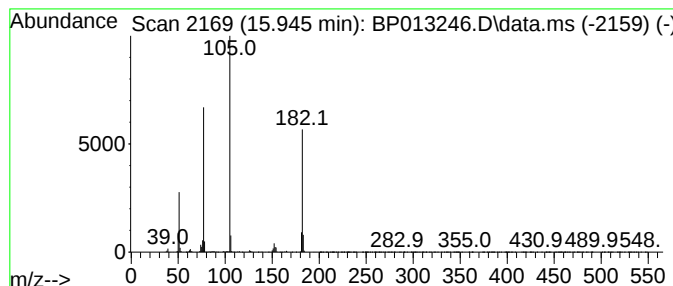
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Benzophenone Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.945	11.14 ng/ul	3346450	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			Azobenzene	182	C12H10N2	000103-33-3	64



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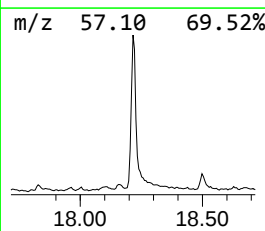
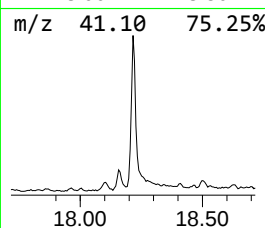
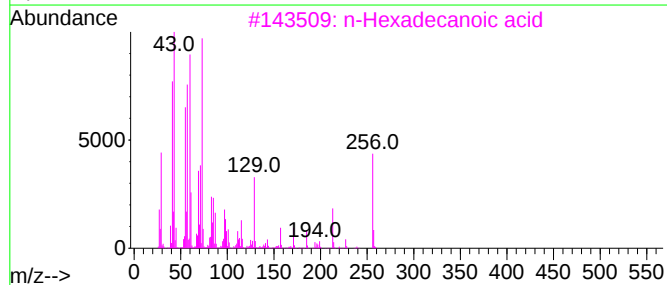
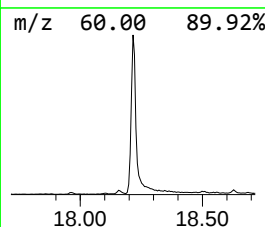
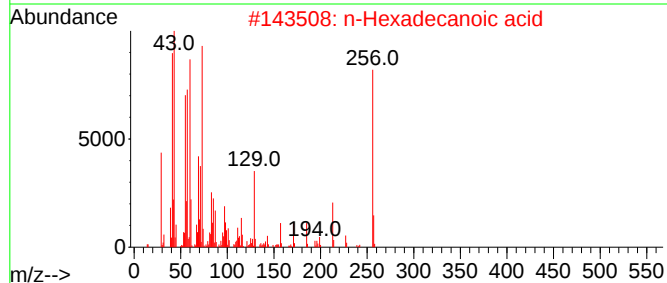
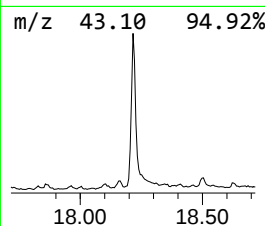
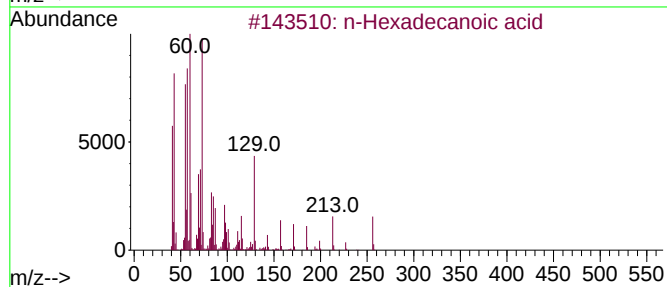
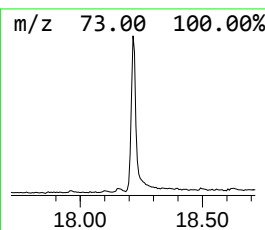
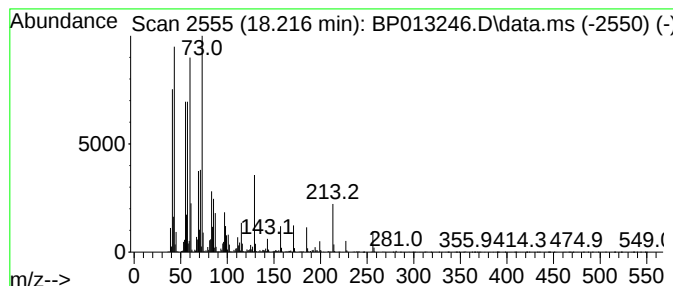
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 n-Hexadecanoic acid Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.216	8.49 ng/ul	2600500	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	99
3			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
4			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	97
5			Tetradecanoic acid	228	C14H28O2	000544-63-8	93



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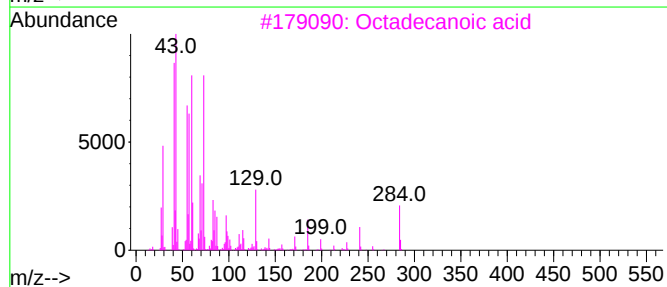
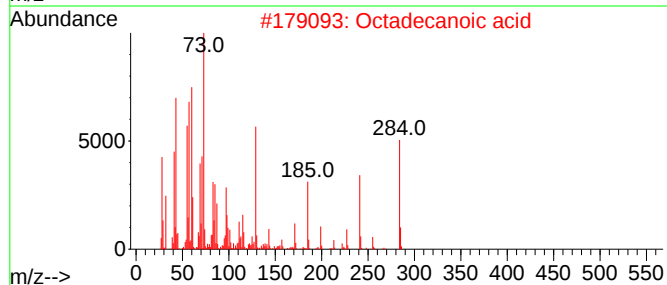
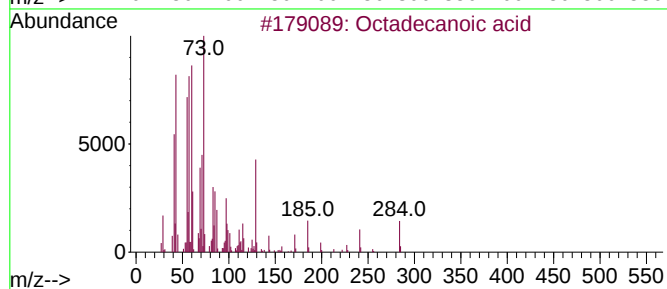
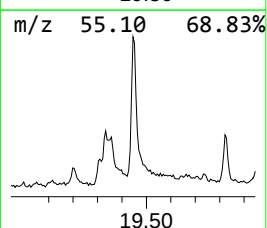
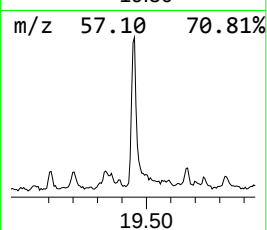
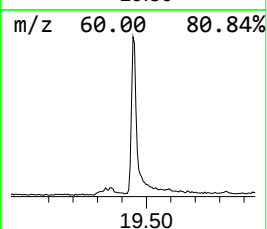
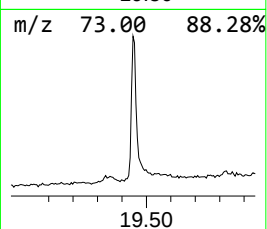
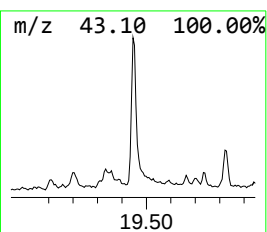
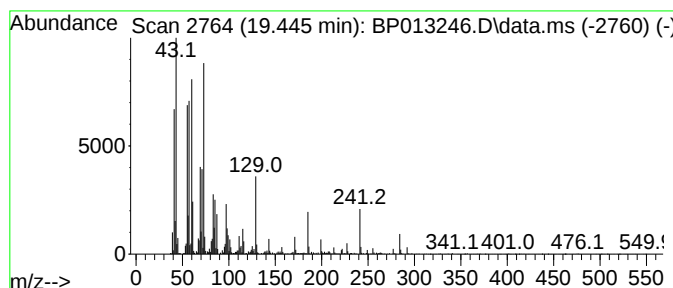
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 Octadecanoic acid Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.445	4.73 ng/ul	1456540	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	98
3			Octadecanoic acid	284	C18H36O2	000057-11-4	98
4			Octadecanoic acid	284	C18H36O2	000057-11-4	94
5			Octadecanoic acid	284	C18H36O2	000057-11-4	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122222\
Data File : BP013246.D
Acq On : 22 Dec 2022 15:09
Operator : CG/JU
Sample : N6015-05
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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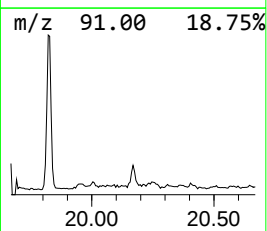
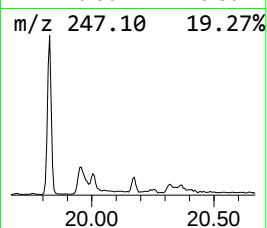
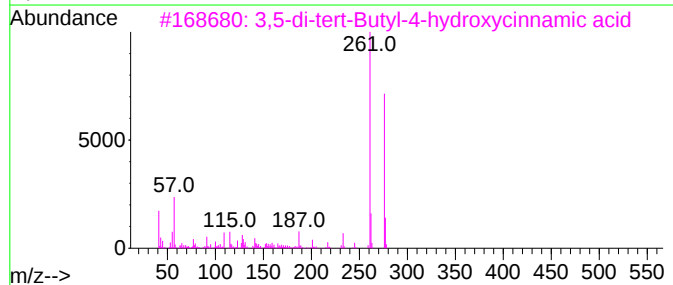
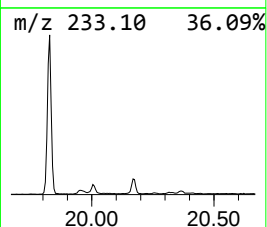
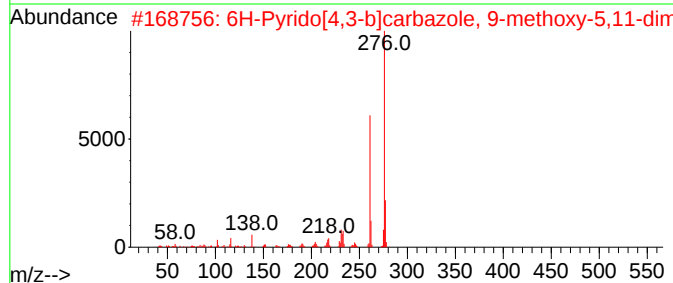
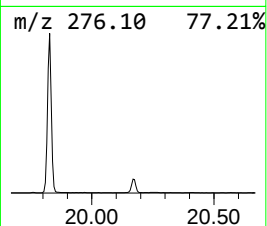
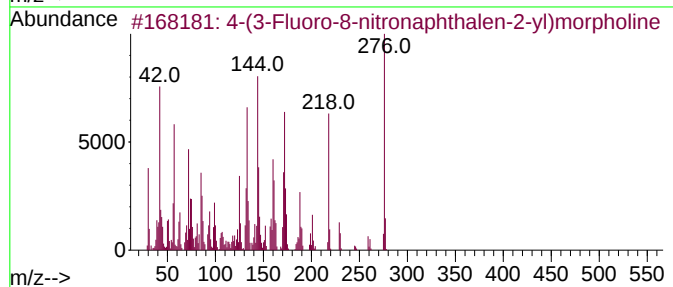
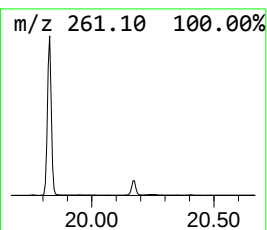
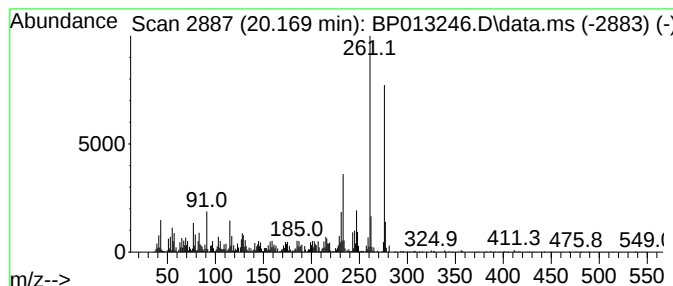
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 6H-Pyrido[4,3-b]carbazole, ... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.169	2.35 ng/ul	724355	Chrysene-d12	21.433

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			4-(3-Fluoro-8-nitronaphthalen-2-yl)morpholine	276	C14H13FN2O3	1000409-37-1	93
2			6H-Pyrido[4,3-b]carbazole, 9-methoxy-5,11-dimethyl-	276	C18H16N2O	010371-86-5	90
3			3,5-di-tert-Butyl-4-hydroxycinnamic acid	276	C17H24O3	022014-01-3	89
4			3-Methoxy-5-pentyl-2-prenylphenol	276	C18H28O2	083036-53-7	81
5			5H-Furo[3,2-g][1]benzopyran-5-one	276	C14H12O6	000668-10-0	72



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122222\
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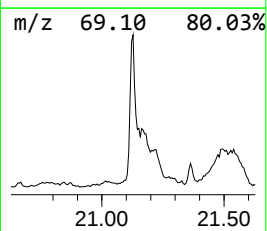
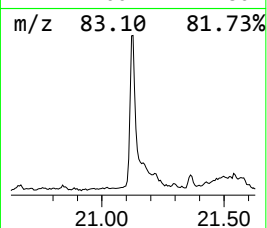
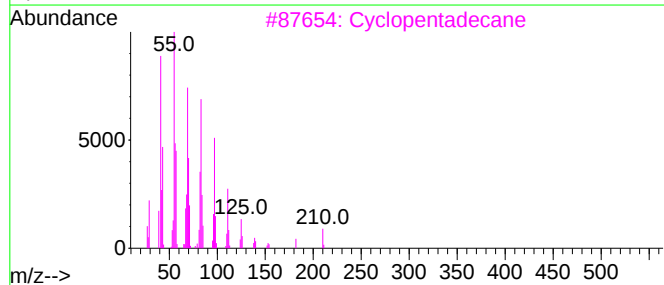
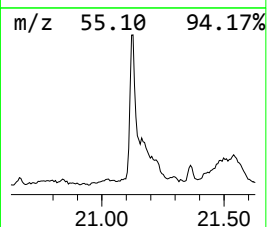
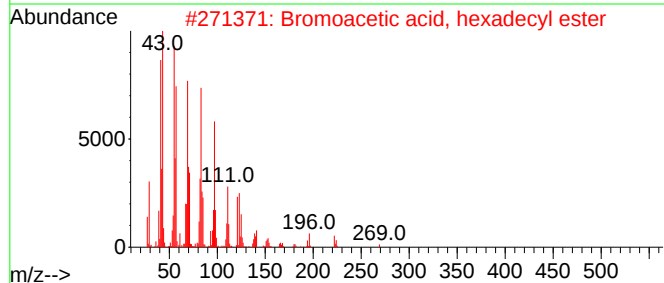
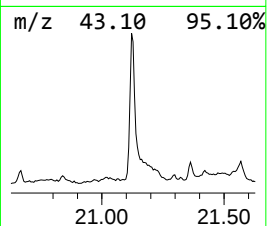
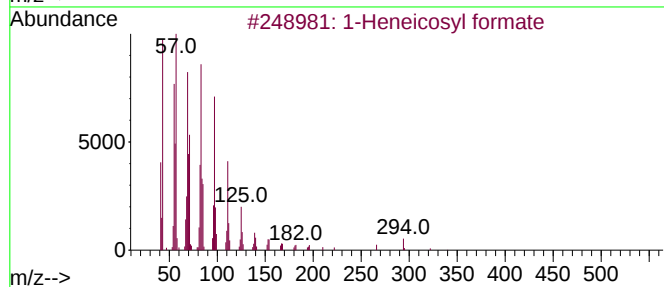
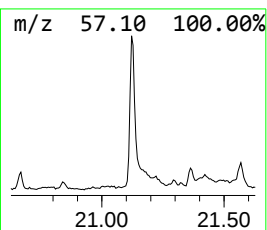
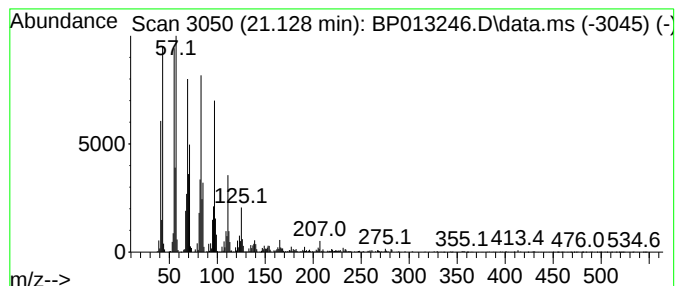
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 1-Heneicosyl formate Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.128	9.16 ng/ul	2822580	Chrysene-d12	21.433

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Heneicosyl formate	340	C22H44O2	077899-03-7	94
2			Bromoacetic acid, hexadecyl ester	362	C18H35BrO2	005454-48-8	94
3			Cyclopentadecane	210	C15H30	000295-48-7	93
4			Heptafluorobutyric acid, pentade...	424	C19H31F7O2	959261-23-5	91
5			Heptacos-1-ene	378	C27H54	015306-27-1	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122222\
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Sample : N6015-05
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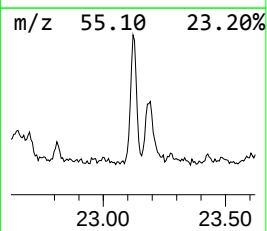
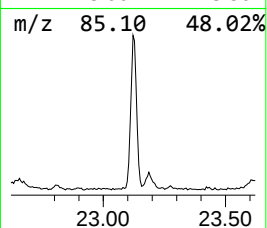
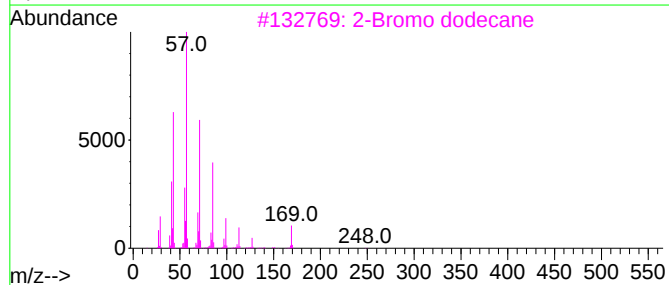
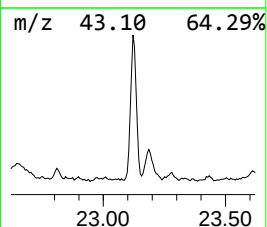
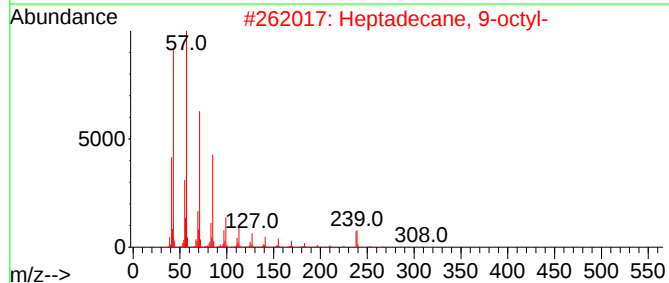
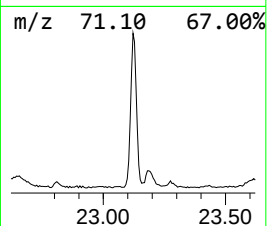
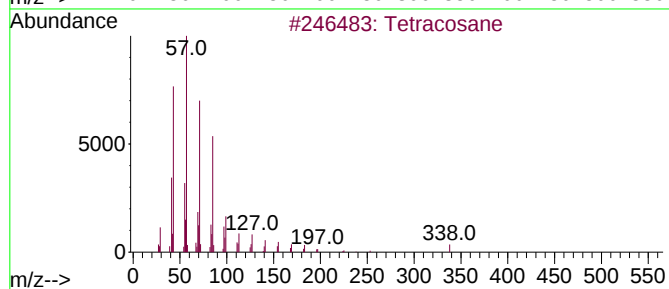
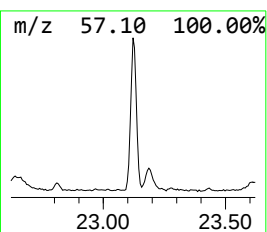
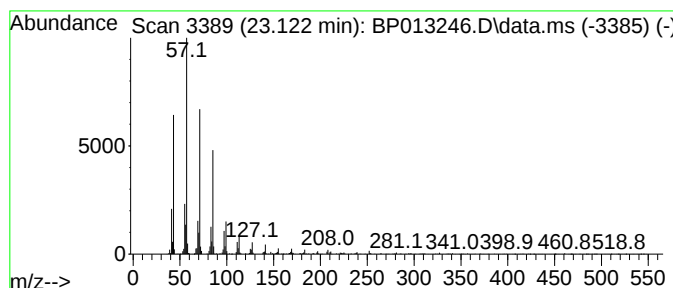
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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.	
23.122	4.17 ng/ul	1370620	Perylene-d12	23.910	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tetracosane	338	C24H50	000646-31-1	93
2	Heptadecane, 9-octyl-	352	C25H52	007225-64-1	93
3	2-Bromo dodecane	248	C12H25Br	013187-99-0	91
4	Heneicosane	296	C21H44	000629-94-7	91
5	Octacosane	394	C28H58	000630-02-4	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122222\
Data File : BP013246.D
Acq On : 22 Dec 2022 15:09
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Sample : N6015-05
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Instrument :
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ClientSampleId :
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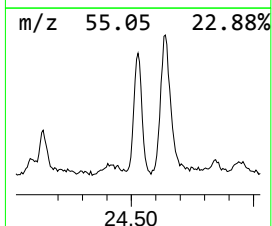
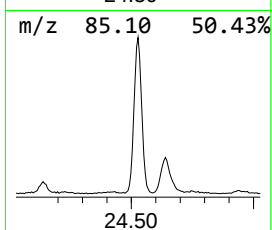
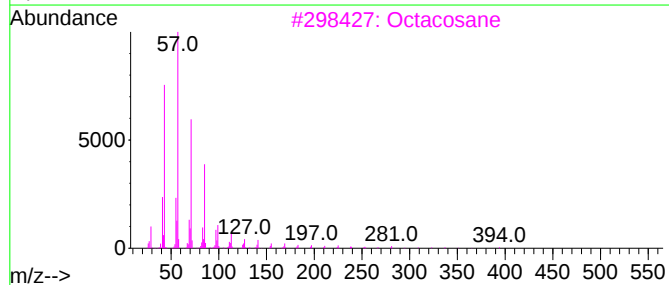
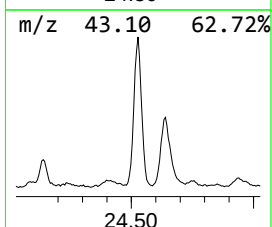
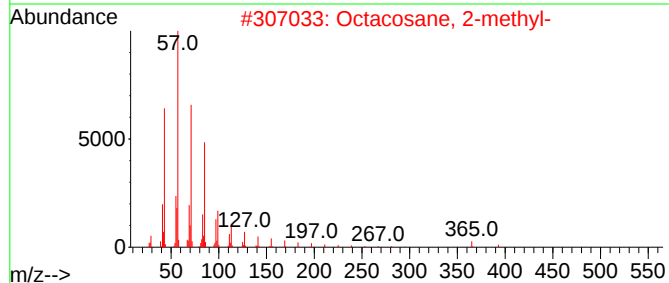
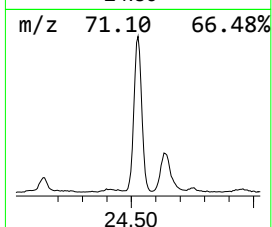
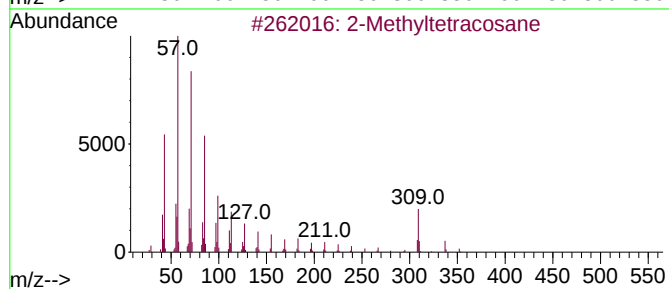
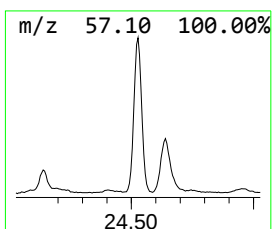
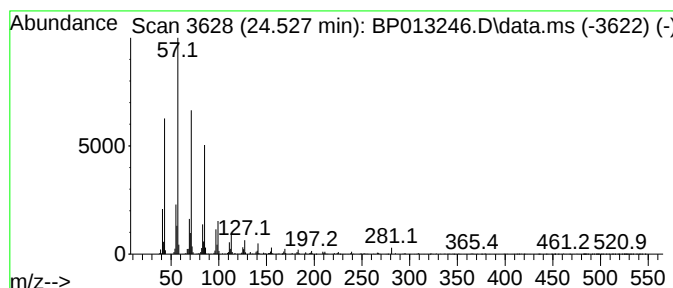
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 (DEL) Alkane: Straight-Chai... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.527	9.89 ng/ul	3247480	Perylene-d12	23.910

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		2-Methyltetracosane	352	C25H52	001560-78-7	94
2		Octacosane, 2-methyl-	408	C29H60	001560-98-1	91
3		Octacosane	394	C28H58	000630-02-4	91
4		Heneicosane	296	C21H44	000629-94-7	91
5		Tetracosane	338	C24H50	000646-31-1	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122222\
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Misc :
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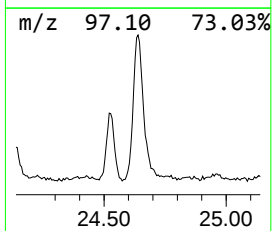
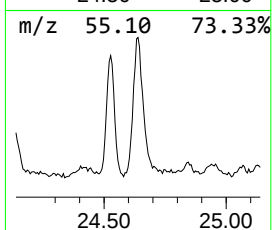
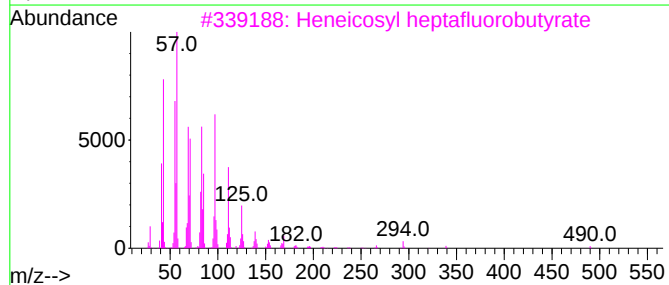
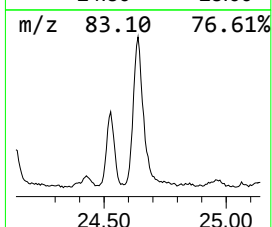
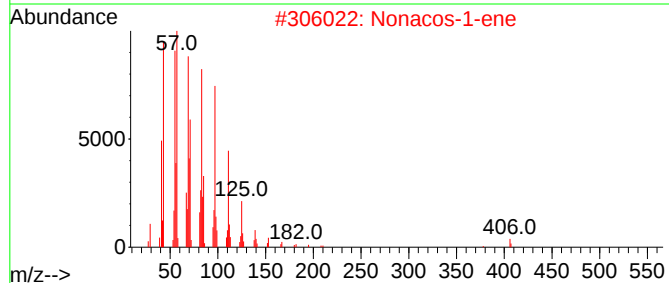
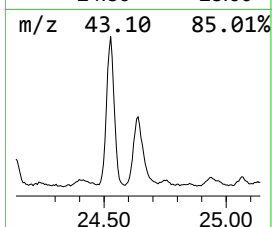
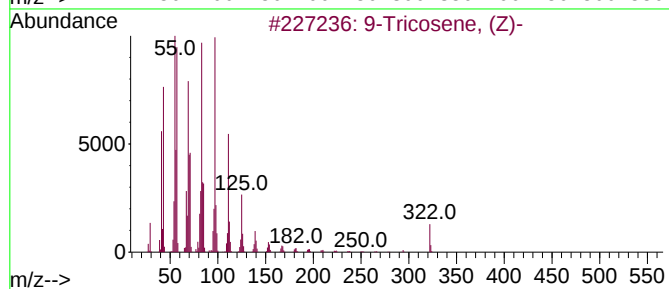
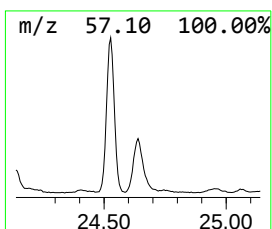
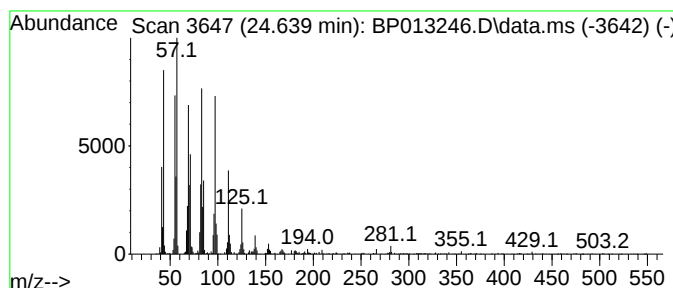
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 (DEL) Alkane: Straight-Chai... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD		R.T.	
24.639	7.96 ng/ul	2614160	Perylene-d12		23.910	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	9-Tricosene, (Z)-		322	C23H46	027519-02-4	96
2	Nonacos-1-ene		406	C29H58	018835-35-3	94
3	Heneicosyl heptafluorobutyrate		508	C25H43F7O2	1000351-83-8	94
4	1-Docosene		308	C22H44	001599-67-3	91
5	Heptacos-1-ene		378	C27H54	015306-27-1	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122222\
Data File : BP013246.D
Acq On : 22 Dec 2022 15:09
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Instrument :
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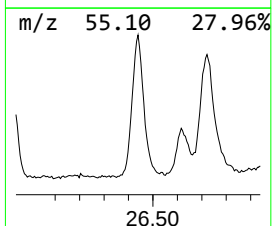
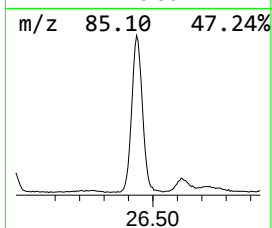
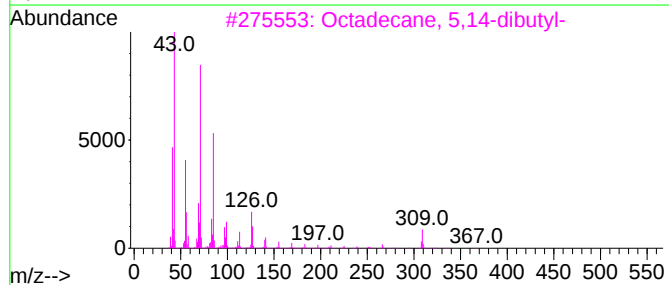
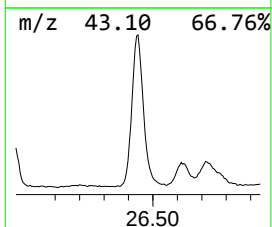
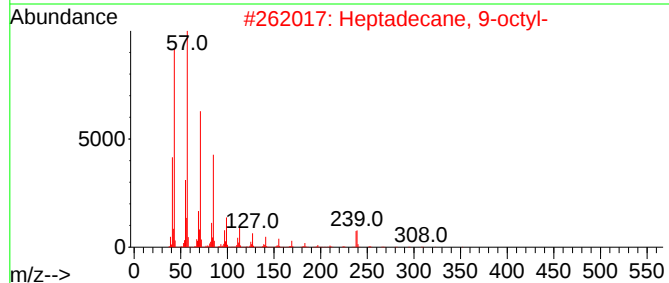
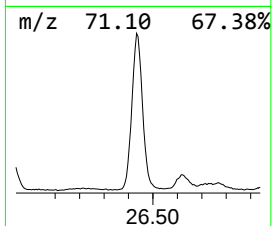
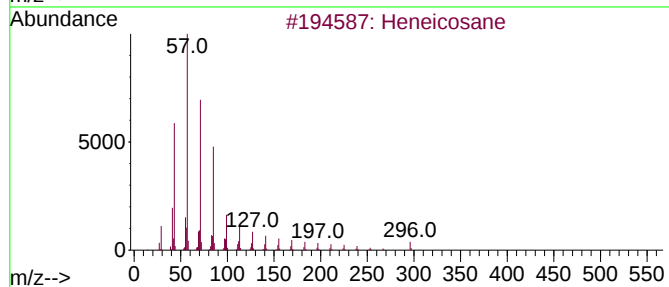
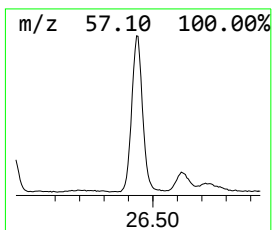
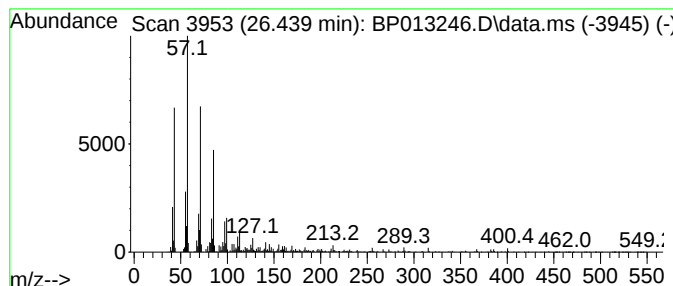
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 (DEL) Alkane: Straight-Chai... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
26.439	23.38 ng/ul	7677360	Perylene-d12	23.910

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heneicosane	296	C21H44	000629-94-7	97
2			Heptadecane, 9-octyl-	352	C25H52	007225-64-1	94
3			Octadecane, 5,14-dibutyl-	366	C26H54	055282-13-8	93
4			2-Methylpentacosane	366	C26H54	000629-87-8	93
5			Docosane, 11-butyl-	366	C26H54	013475-76-8	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122222\
Data File : BP013246.D
Acq On : 22 Dec 2022 15:09
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Sample : N6015-05
Misc :
ALS Vial : 13 Sample Multiplier: 1

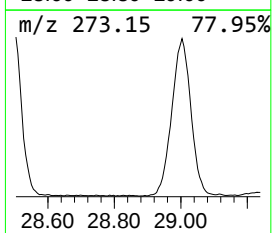
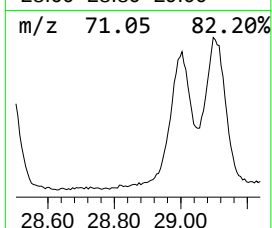
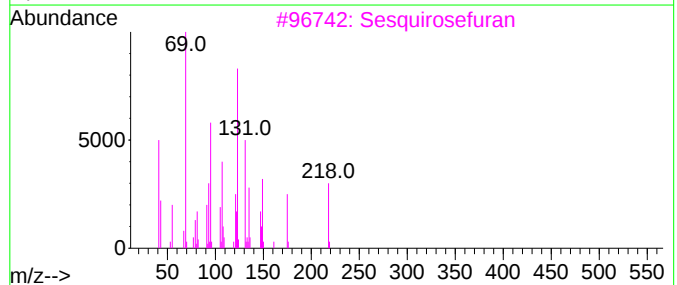
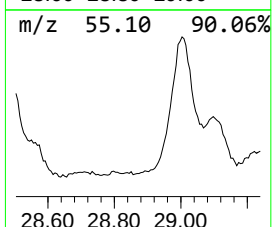
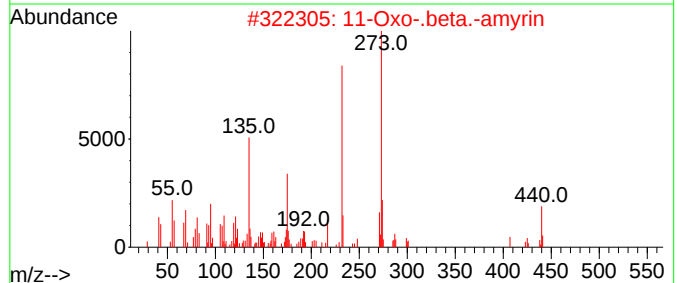
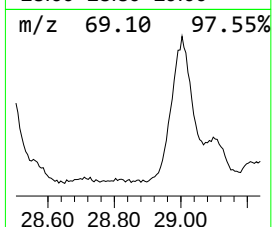
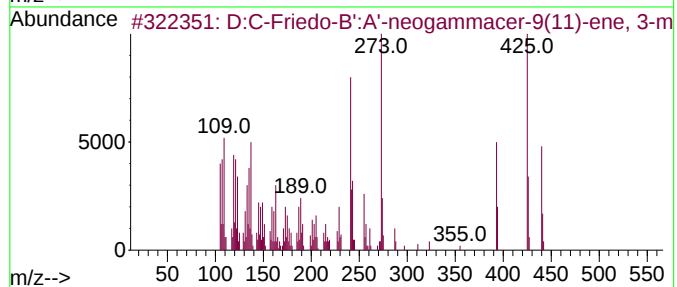
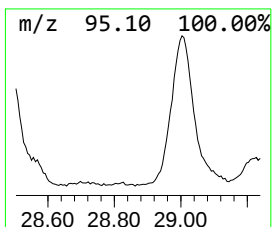
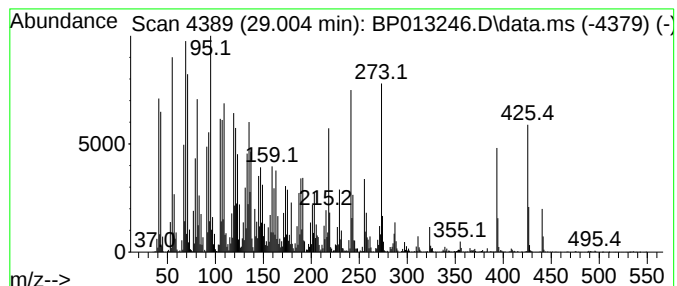
Instrument :
BNA_P
ClientSampleId :
DBXU1

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 11 D:C-Friedo-B':A'-neogammace... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.	
29.004	35.04 ng/ul	11506100	Perylene-d12	23.910	
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	11-Oxo-.beta.-amyrin	440	C30H48O2	038242-02-3	46
3	Sesquirosefuran	218	C15H22O	039007-93-7	35
4	9-Thiabicyclo[3.3.1]nonane, 2,6-...	274	C14H18N4S	1000141-52-1	27
5	Lupeol	426	C30H50O	000545-47-1	20



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122222\
Data File : BP013246.D
Acq On : 22 Dec 2022 15:09
Operator : CG/JU
Sample : N6015-05
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DBXU1

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	11.1	ng/ul	3346450	3	14.587	6009750	20.0
n-Hexadecanoic ...	18.216	8.5	ng/ul	2600500	4	17.345	6127000	20.0
Octadecanoic acid	19.445	4.7	ng/ul	1456540	5	21.433	6162010	20.0
4-(3-Fluoro-8-n...	19.828	16.7	ng/ul	5135340	5	21.433	6162010	20.0
6H-Pyrido[4,3-b...	20.169	2.4	ng/ul	724355	5	21.433	6162010	20.0
1-Heneicosyl fo...	21.128	9.2	ng/ul	2822580	5	21.433	6162010	20.0
(DEL) Alkane: S...	23.122	4.2	ng/ul	1370620	6	23.910	6566760	20.0
(DEL) Alkane: S...	24.527	9.9	ng/ul	3247480	6	23.910	6566760	20.0
(DEL) Alkane: S...	24.639	8.0	ng/ul	2614160	6	23.910	6566760	20.0
(DEL) Alkane: S...	26.439	23.4	ng/ul	7677360	6	23.910	6566760	20.0
D:C-Friedo-B':A...	29.004	35.0	ng/ul	11506100	6	23.910	6566760	20.0