

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122222\
 Data File : BP013245.D
 Acq On : 22 Dec 2022 14:34
 Operator : CG/JU
 Sample : N6015-04
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 DBXU0

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: BP013245.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	34	rVB	140162	190826	1.83%	0.115%
2	3.770	97	99	114	rBV	288429	487724	4.67%	0.293%
3	3.987	134	136	142	rVB3	19435	30787	0.29%	0.019%
4	4.093	149	154	159	rVB	52447	69565	0.67%	0.042%
5	5.040	310	315	320	rVB3	27512	44498	0.43%	0.027%
6	5.270	352	354	360	rVB4	12851	19798	0.19%	0.012%
7	7.111	659	667	679	rBV	2061605	3323717	31.80%	1.999%
8	7.275	689	695	707	rVB	1868796	2886069	27.62%	1.735%
9	7.475	720	729	739	rBV	2249767	3553753	34.01%	2.137%
10	7.946	801	809	816	rBV	2194333	3535253	33.83%	2.126%
11	8.005	816	819	826	rVB	18020	27363	0.26%	0.016%
12	8.658	923	930	943	rBV	2263074	3650731	34.93%	2.195%
13	9.116	1001	1008	1015	rBV	2238986	3616370	34.60%	2.175%
14	9.275	1030	1035	1041	rVB7	25799	41030	0.39%	0.025%
15	9.522	1070	1077	1083	rVB	67866	111735	1.07%	0.067%
16	9.710	1104	1109	1114	rVB3	11354	21338	0.20%	0.013%
17	9.840	1122	1131	1143	rBV	1849454	3003433	28.74%	1.806%
18	10.069	1164	1170	1174	rVB8	12283	22490	0.22%	0.014%
19	10.263	1199	1203	1209	rVB7	11937	19922	0.19%	0.012%
20	10.381	1215	1223	1232	rBV	2850647	4679058	44.77%	2.814%
21	10.763	1278	1288	1295	rBV	3170751	5280474	50.53%	3.175%
22	10.899	1303	1311	1322	rBV	1544786	2717316	26.00%	1.634%
23	11.299	1376	1379	1385	rVB7	10157	18380	0.18%	0.011%
24	11.552	1417	1422	1425	rBV7	11554	21950	0.21%	0.013%
25	11.669	1436	1442	1445	rBV8	13592	25869	0.25%	0.016%
26	11.781	1456	1461	1467	rVB5	18263	36280	0.35%	0.022%
27	12.016	1497	1501	1506	rVB4	16877	28412	0.27%	0.017%
28	12.075	1506	1511	1517	rBV5	14036	28667	0.27%	0.017%
29	12.169	1522	1527	1535	rVV	46411	77001	0.74%	0.046%
30	12.246	1538	1540	1546	rVV5	12843	20955	0.20%	0.013%
31	12.346	1552	1557	1564	rVV2	53906	97973	0.94%	0.059%
32	12.416	1565	1569	1578	rVB4	21413	42093	0.40%	0.025%
33	12.557	1585	1593	1600	rBV4	7553	19390	0.19%	0.012%
34	12.634	1600	1606	1613	rBV5	20546	40005	0.38%	0.024%
35	12.799	1629	1634	1638	rBV6	11528	17987	0.17%	0.011%
36	12.981	1661	1665	1669	rVB4	14088	22059	0.21%	0.013%

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Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
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37	13.116	1684	1688	1697	rVV4	25131	47128	0.45%	0.028%
38	13.328	1720	1724	1728	rBV7	13899	21084	0.20%	0.013%
39	13.404	1728	1737	1745	rVV	96075	155741	1.49%	0.094%
40	13.569	1763	1765	1772	rVB7	18295	31167	0.30%	0.019%
41	13.769	1792	1799	1802	rBV9	11145	21745	0.21%	0.013%
42	13.916	1817	1824	1830	rBV8	41232	86822	0.83%	0.052%
43	13.998	1830	1838	1846	rVV	4746937	6555456	62.73%	3.942%
44	14.057	1846	1848	1854	rVV4	30280	40371	0.39%	0.024%
45	14.110	1854	1857	1861	rVB6	17147	26989	0.26%	0.016%
46	14.181	1865	1869	1872	rVB	18890	25074	0.24%	0.015%
47	14.287	1880	1887	1897	rBV	5823663	8362297	80.02%	5.029%
48	14.475	1914	1919	1927	rVB	61312	100478	0.96%	0.060%
49	14.593	1932	1939	1948	rVB	4815273	6998696	66.97%	4.209%
50	14.787	1966	1972	1991	rBV	1044391	1863102	17.83%	1.120%
51	14.945	1996	1999	2003	rVV6	21608	39147	0.37%	0.024%
52	14.993	2003	2007	2014	rVV5	49600	85841	0.82%	0.052%
53	15.381	2068	2073	2075	rVV6	17041	31893	0.31%	0.019%
54	15.428	2077	2081	2087	rVB3	28245	47806	0.46%	0.029%
55	15.587	2101	2108	2120	rBV	7228072	10340912	98.95%	6.218%
56	15.704	2122	2128	2136	rBV	1693173	2233801	21.37%	1.343%
57	15.828	2145	2149	2154	rVB3	39604	58850	0.56%	0.035%
58	15.951	2164	2170	2184	rVB	2919171	3837395	36.72%	2.308%
59	16.216	2212	2215	2220	rBV7	19900	29454	0.28%	0.018%
60	16.310	2225	2231	2236	rVV6	42853	99106	0.95%	0.060%
61	16.451	2249	2255	2262	rBV10	19214	62106	0.59%	0.037%
62	16.522	2262	2267	2274	rVV2	65719	101637	0.97%	0.061%
63	16.734	2298	2303	2308	rBV6	45117	76012	0.73%	0.046%
64	17.092	2360	2364	2368	rBV	29808	44439	0.43%	0.027%
65	17.134	2368	2371	2374	rBV3	14290	20074	0.19%	0.012%
66	17.228	2383	2387	2391	rBV3	81459	111088	1.06%	0.067%
67	17.298	2395	2399	2401	rBV3	39663	47763	0.46%	0.029%
68	17.351	2401	2408	2418	rBV	5537346	8461988	80.97%	5.088%
69	17.445	2418	2424	2432	rVB2	6875856	9846633	94.22%	5.921%
70	17.622	2450	2454	2461	rBV7	51825	90726	0.87%	0.055%
71	17.704	2465	2468	2470	rBV3	31417	39313	0.38%	0.024%
72	17.834	2487	2490	2492	rBV3	26393	27232	0.26%	0.016%
73	18.004	2516	2519	2522	rVB4	32947	37127	0.36%	0.022%
74	18.104	2533	2536	2541	rVB4	65648	88503	0.85%	0.053%
75	18.163	2542	2546	2551	rBV	164737	250953	2.40%	0.151%
76	18.222	2551	2556	2567	rBV	1608952	2306654	22.07%	1.387%

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77	18.410	2582	2588	2593	rVB	615821	811136	7.76%	0.488%
78	18.469	2594	2598	2601	rBV	253060	320275	3.06%	0.193%
79	18.498	2601	2603	2608	rVB3	63522	83732	0.80%	0.050%
80	18.628	2621	2625	2632	rBV	442076	568904	5.44%	0.342%
81	18.722	2637	2641	2647	rVB	291750	360875	3.45%	0.217%
82	18.786	2647	2652	2657	rBV	1362757	1617544	15.48%	0.973%
83	19.122	2704	2709	2713	rBV2	514145	664542	6.36%	0.400%
84	19.251	2729	2731	2735	rVB2	101795	116289	1.11%	0.070%
85	19.333	2738	2745	2746	rBV3	206882	356863	3.41%	0.215%
86	19.451	2761	2765	2772	rBV	589462	832040	7.96%	0.500%
87	19.681	2798	2804	2816	rBV	7693713	10223152	97.82%	6.148%
88	19.828	2824	2829	2837	rBV	8803647	10450639	100.00%	6.284%
89	19.951	2846	2850	2857	rBV	292617	551975	5.28%	0.332%
90	20.175	2883	2888	2894	rVB2	1391712	1818727	17.40%	1.094%
91	20.410	2925	2928	2932	rBV4	135311	154870	1.48%	0.093%
92	20.669	2970	2972	2976	rVB	214379	209602	2.01%	0.126%
93	21.128	3046	3050	3061	rBV	1642994	2299914	22.01%	1.383%
94	21.433	3097	3102	3108	rBV	4469527	5957296	57.00%	3.582%
95	23.127	3386	3390	3396	rBV2	574303	893865	8.55%	0.538%
96	23.751	3490	3496	3511	rVB2	4089955	9048744	86.59%	5.441%
97	23.910	3517	3523	3536	rVB2	3032856	6293270	60.22%	3.784%
98	24.533	3624	3629	3638	rVB	954229	2000175	19.14%	1.203%
99	26.439	3947	3953	3974	rVB	1365308	4603040	44.05%	2.768%
100	29.009	4382	4390	4402	rVB8	1594593	5476788	52.41%	3.293%

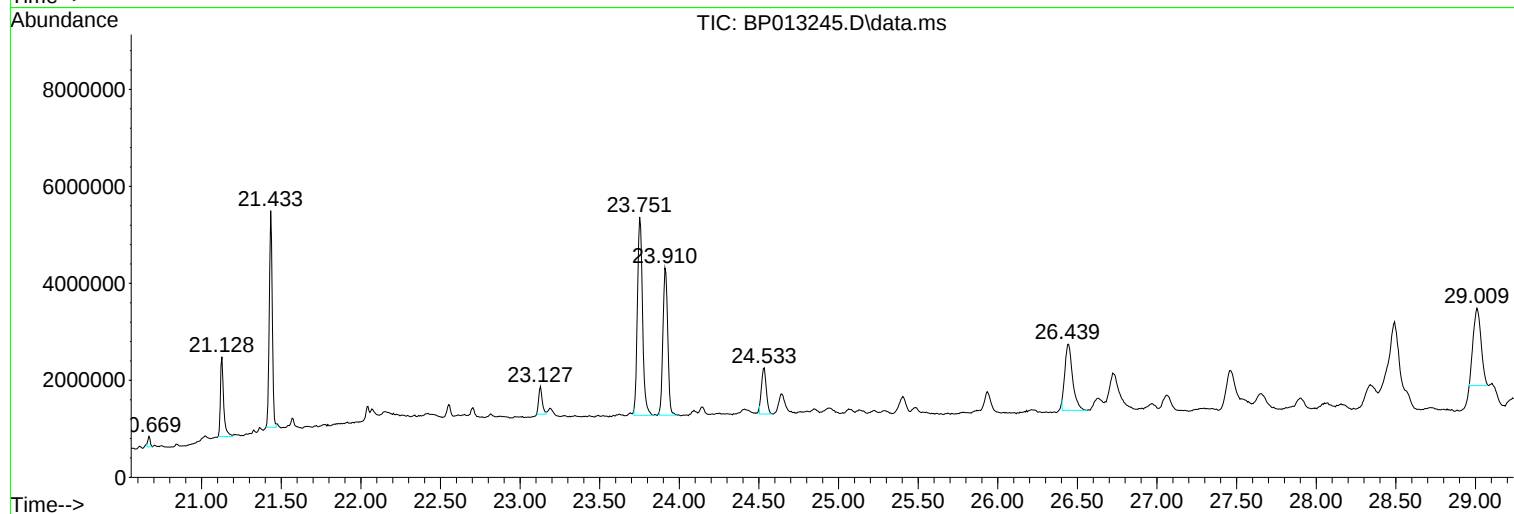
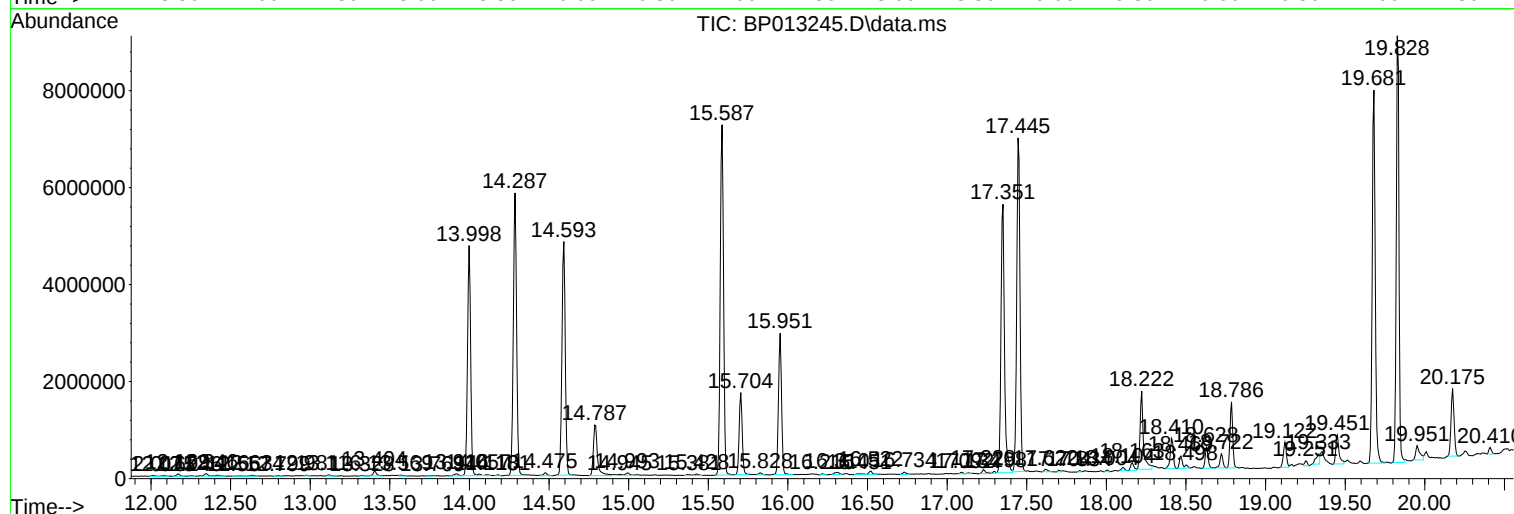
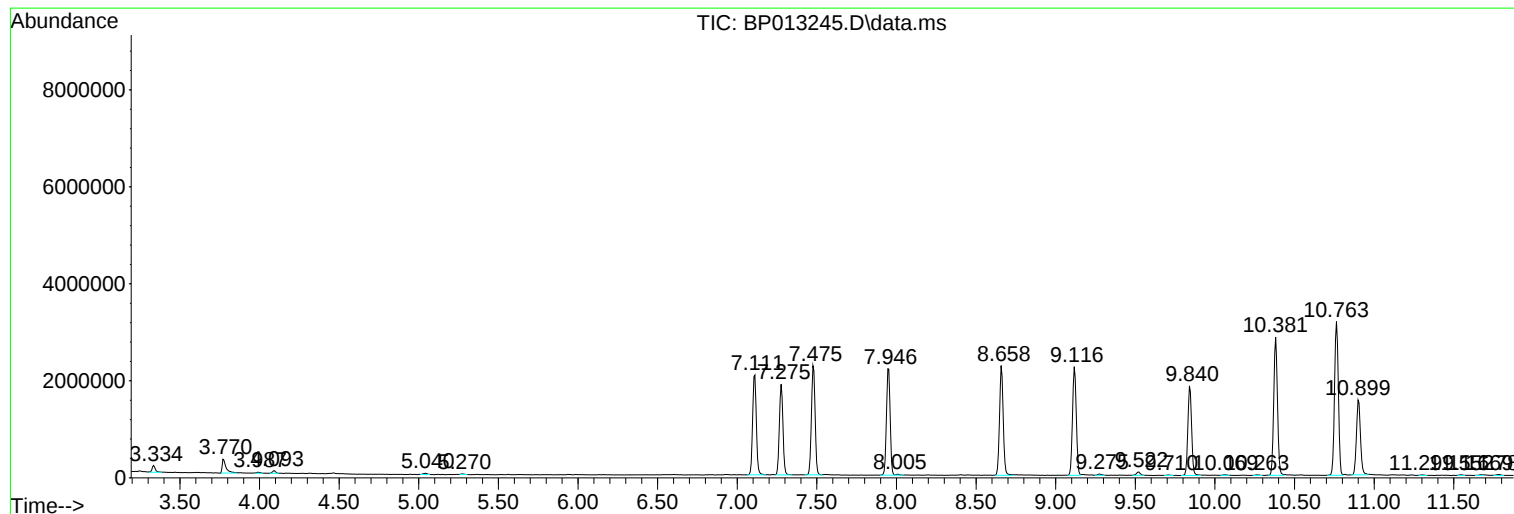
Sum of corrected areas: 166297131

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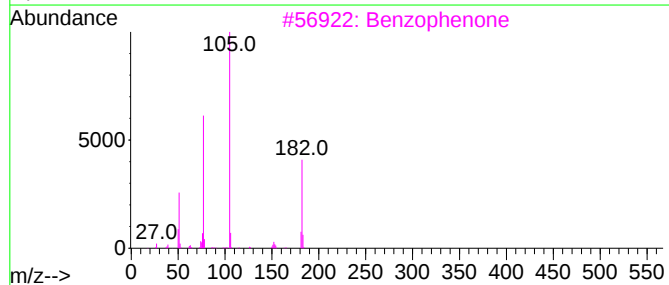
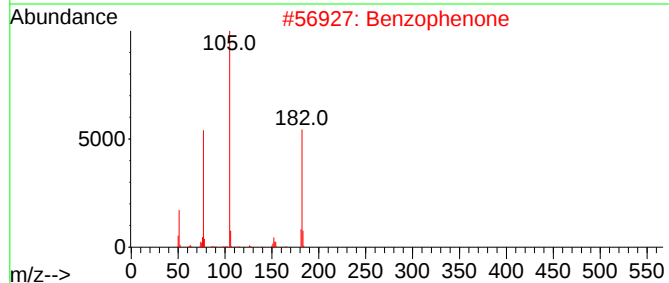
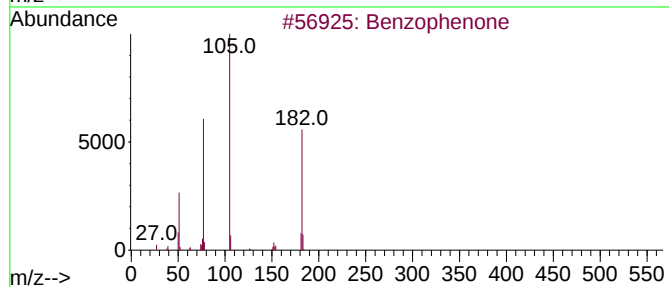
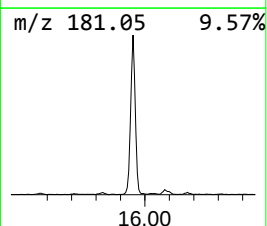
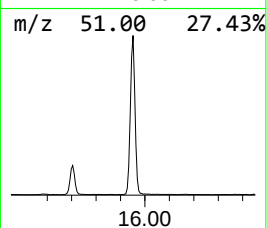
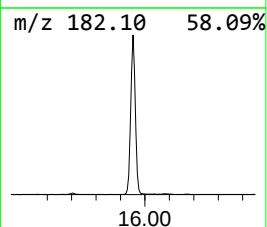
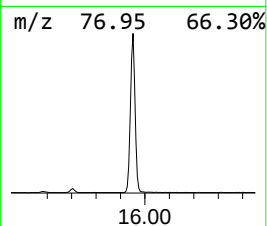
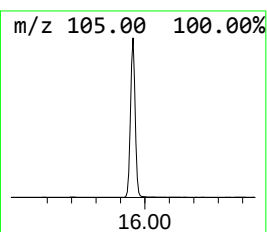
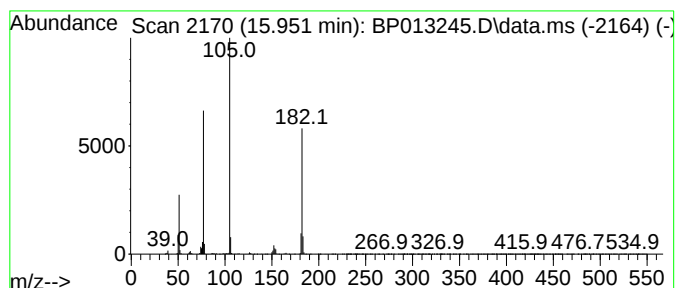
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Benzophenone Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.951	10.97 ng/ul	3837400	Acenaphthene-d10	14.593

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	94
5			Benzophenone	182	C13H10O	000119-61-9	90



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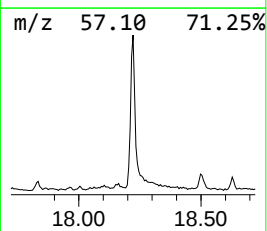
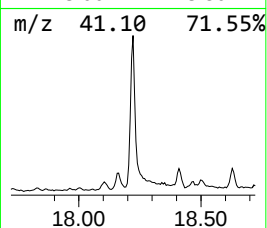
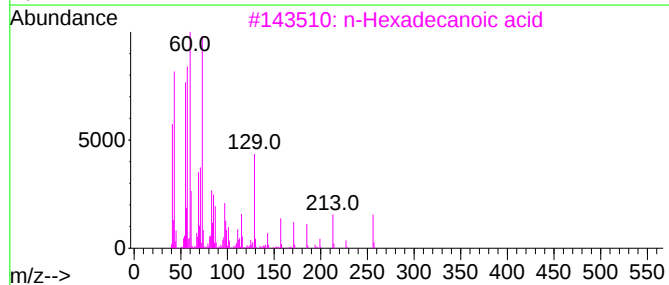
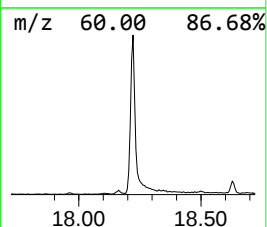
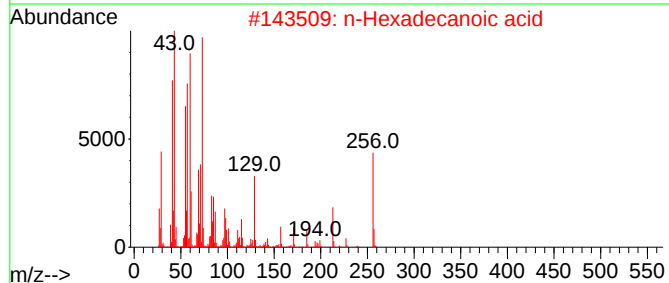
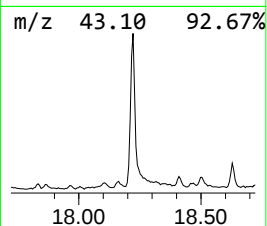
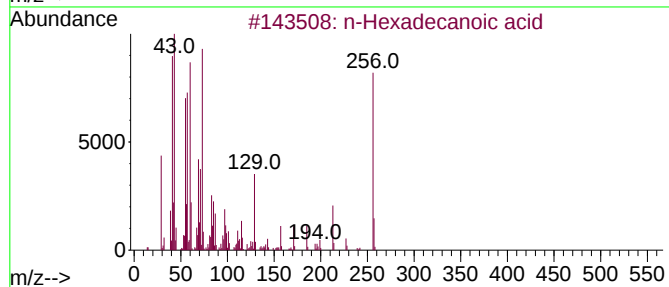
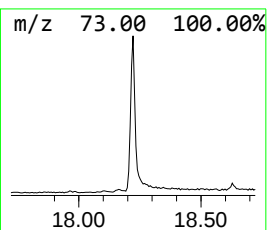
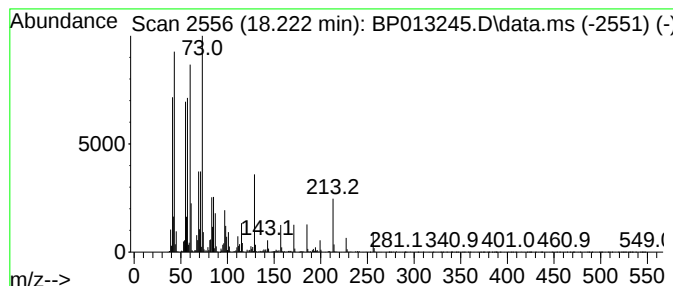
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 n-Hexadecanoic acid Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.222	5.45 ng/ul	2306650	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	98
3			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
4			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	93
5			Pentadecanoic acid	242	C15H30O2	001002-84-2	91



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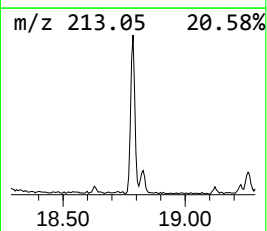
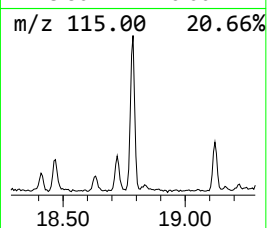
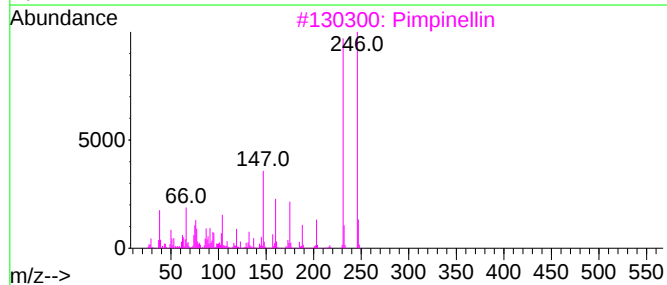
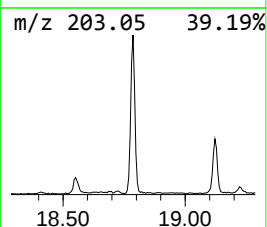
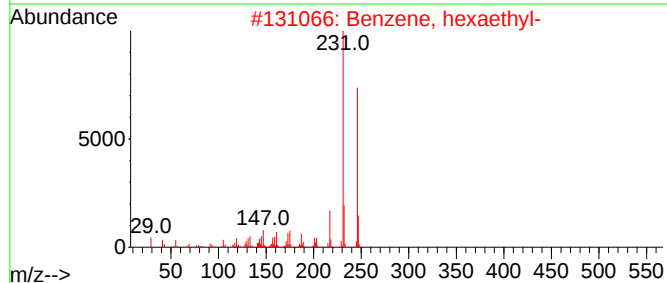
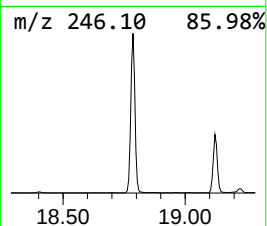
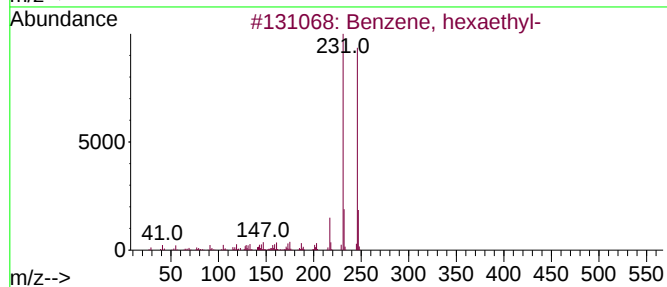
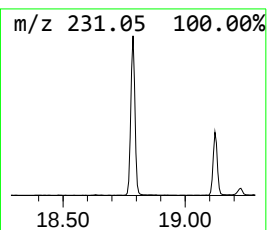
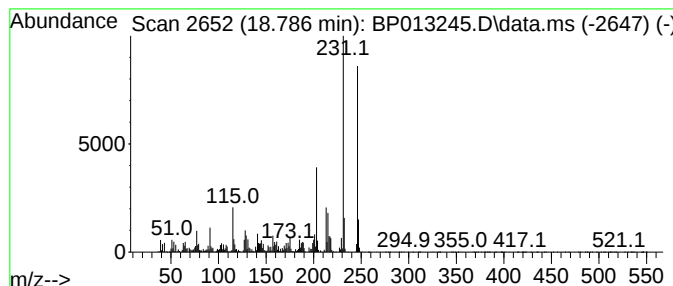
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Quant Title : SVOA CALIBRATION

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

Peak Number 3 Benzene, hexaethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.787	3.82 ng/ul	1617540	Phenanthrene-d10	17.351

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, hexaethyl-	246	C18H30	000604-88-6	86
2			Benzene, hexaethyl-	246	C18H30	000604-88-6	86
3			Pimpinellin	246	C13H10O5	000131-12-4	70
4			7H-Furo[3,2-g][1]benzopyran-7-on...	246	C13H10O5	000482-27-9	62
5			7H-Furo[3,2-g][1]benzopyran-7-on...	246	C13H10O5	000482-27-9	62



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122222\
Data File : BP013245.D
Acq On : 22 Dec 2022 14:34
Operator : CG/JU
Sample : N6015-04
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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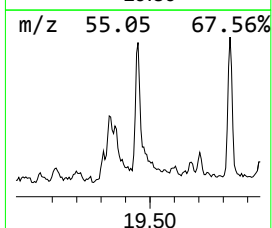
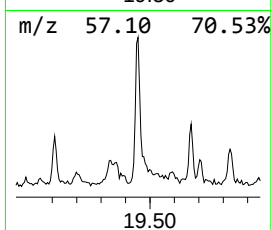
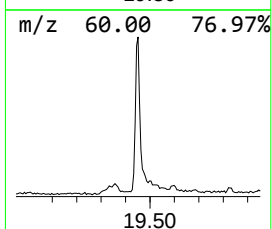
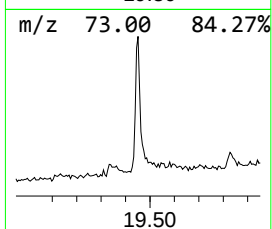
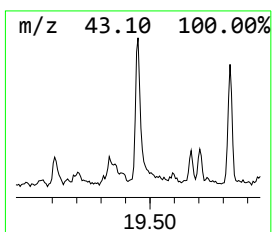
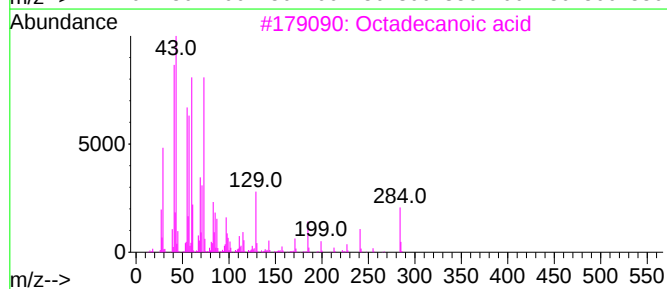
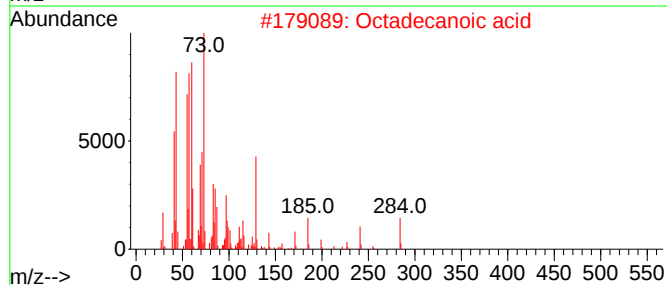
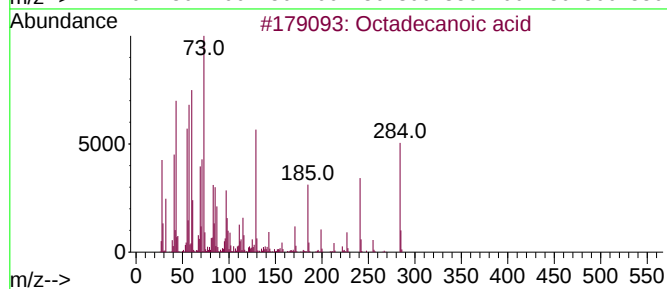
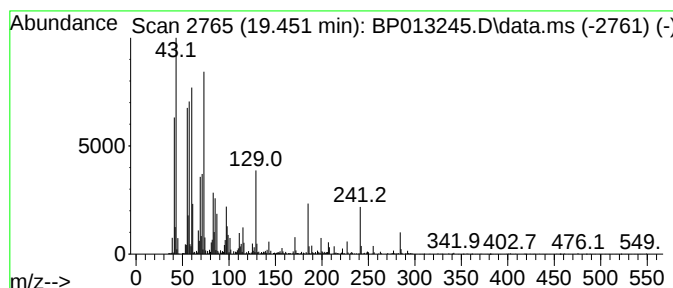
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 Octadecanoic acid Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.451	2.79 ng/ul	832040	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	97
4			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	93
5			Octadecanoic acid	284	C18H36O2	000057-11-4	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122222\
Data File : BP013245.D
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Sample : N6015-04
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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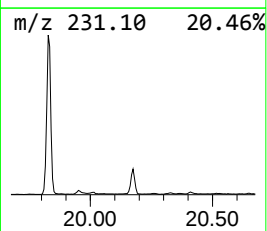
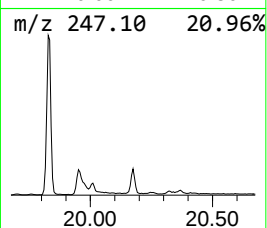
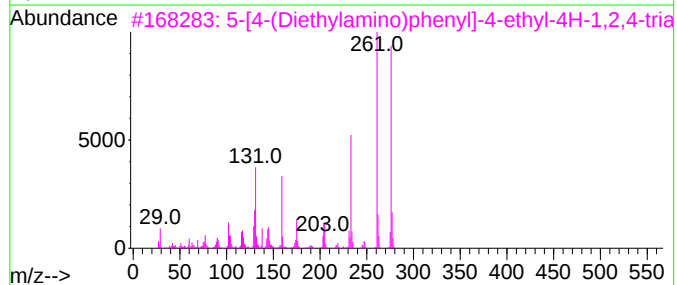
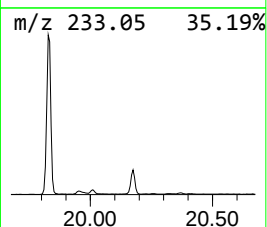
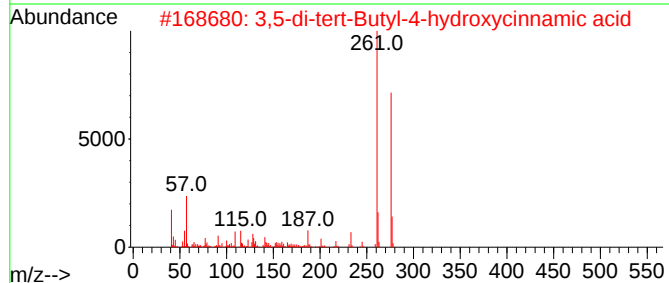
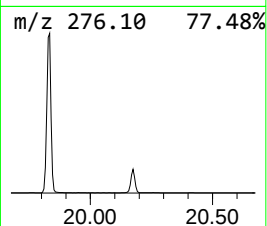
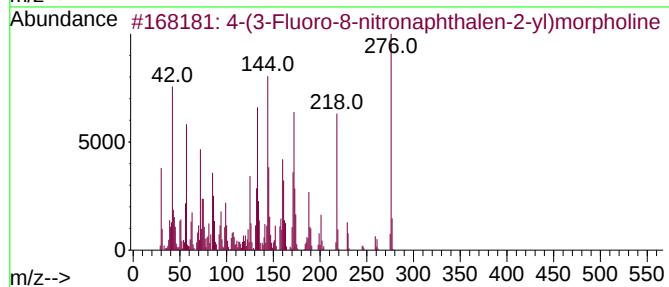
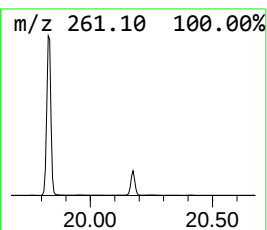
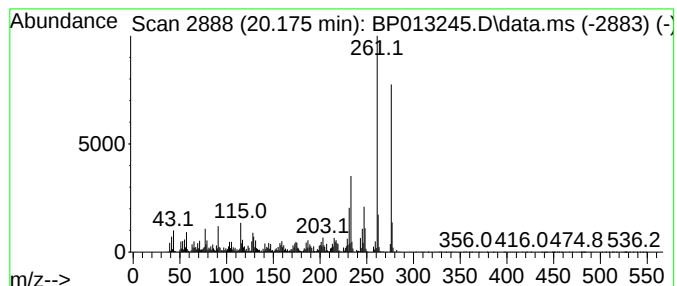
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 3,5-di-tert-Butyl-4-hydroxy... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.175	6.11 ng/ul	1818730	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	91
2			3,5-di-tert-Butyl-4-hydroxycinna...	276	C17H24O3	022014-01-3	70
3			5-[4-(Diethylamino)phenyl]-4-eth...	276	C14H20N4S	1000476-62-2	64
4			6H-Pyrido[4,3-b]carbazole, 9-met...	276	C18H16N2O	010371-86-5	64
5			3,5-di-tert-Butyl-4-hydroxycinna...	276	C17H24O3	022014-01-3	62



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122222\
Data File : BP013245.D
Acq On : 22 Dec 2022 14:34
Operator : CG/JU
Sample : N6015-04
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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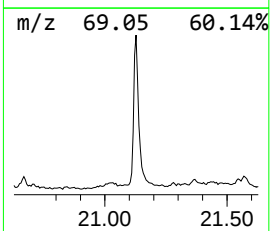
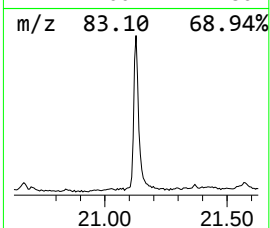
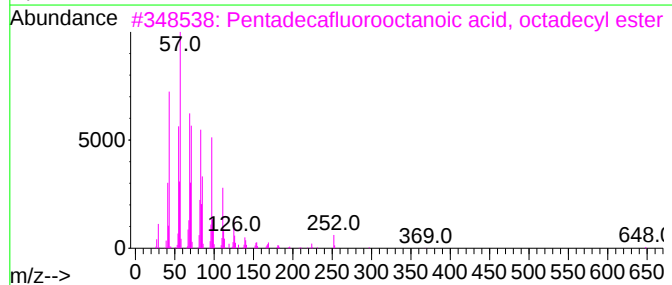
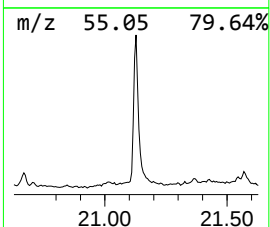
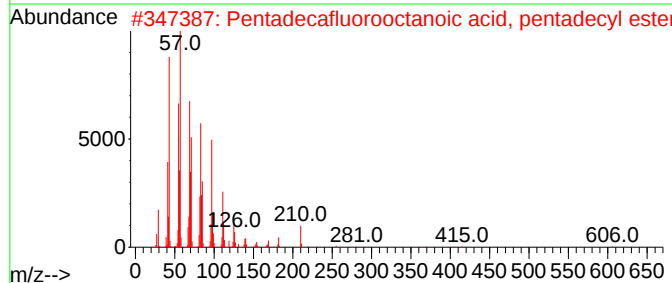
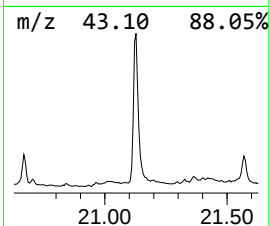
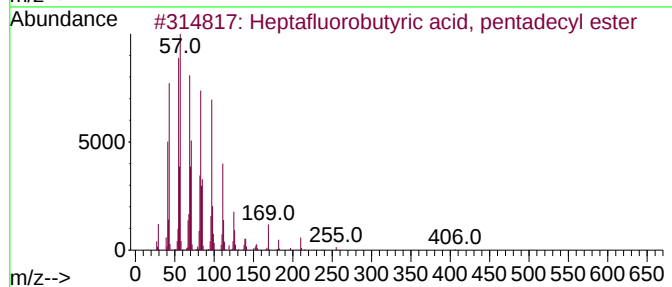
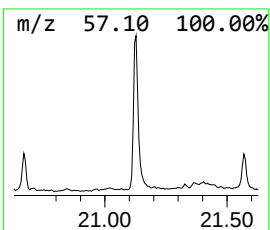
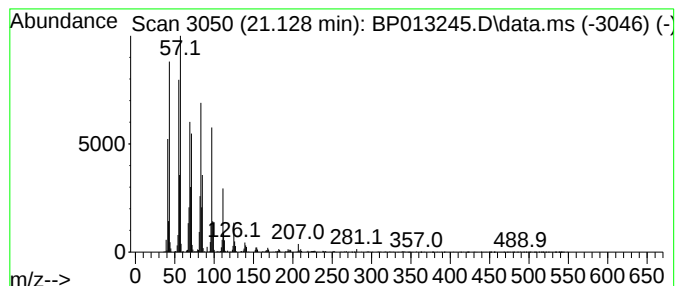
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 Heptafluorobutyric acid, pe... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.128	7.72 ng/ul	2299910	Chrysene-d12	21.433

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heptafluorobutyric acid, pentade...	424	C19H31F7O2	959261-23-5	94
2			Pentadecafluorooctanoic acid, pe...	624	C23H31F15O2	1000406-04-5	94
3			Pentadecafluorooctanoic acid, oc...	666	C26H37F15O2	1000406-04-8	94
4			1-Hexacosene	364	C26H52	018835-33-1	94
5			Heptafluorobutyric acid, n-octad...	466	C22H37F7O2	000400-57-7	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122222\
Data File : BP013245.D
Acq On : 22 Dec 2022 14:34
Operator : CG/JU
Sample : N6015-04
Misc :
ALS Vial : 12 Sample Multiplier: 1

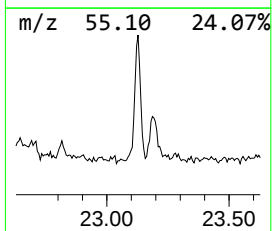
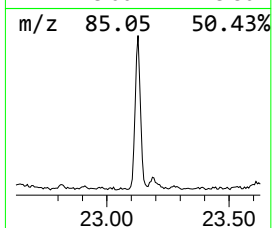
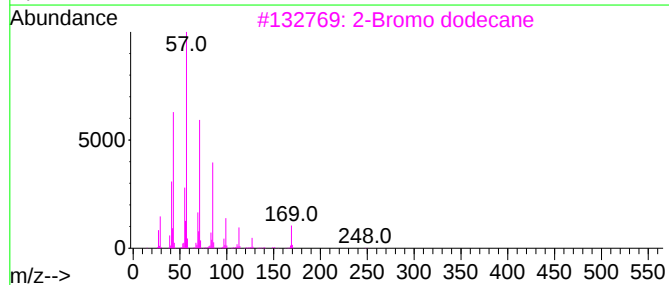
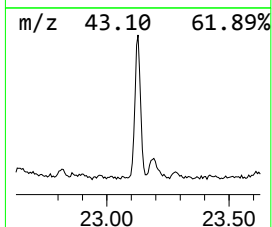
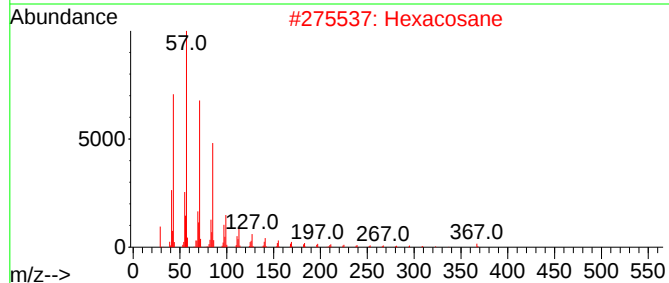
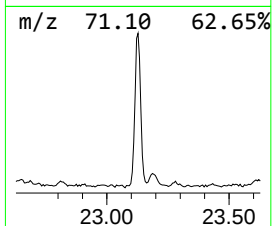
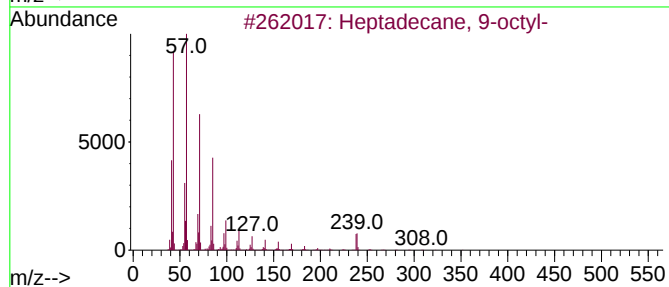
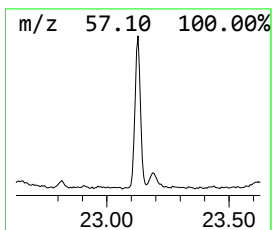
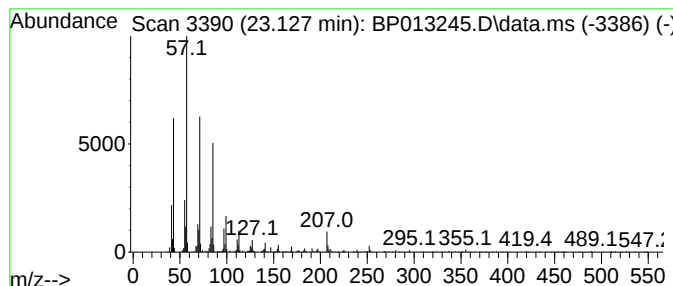
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 (DEL) Alkane: Straight-Chai... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD		R.T.	
23.127	2.84 ng/ul	893865	Perylene-d12		23.910	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Heptadecane, 9-octyl-		352	C25H52	007225-64-1	93
2	Hexacosane		366	C26H54	000630-01-3	93
3	2-Bromo dodecane		248	C12H25Br	013187-99-0	91
4	Tridecane, 1-iodo-		310	C13H27I	035599-77-0	90
5	Heneicosane		296	C21H44	000629-94-7	87



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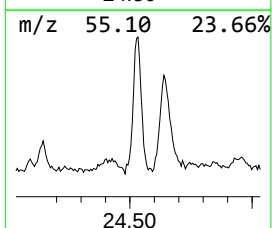
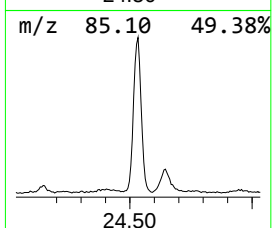
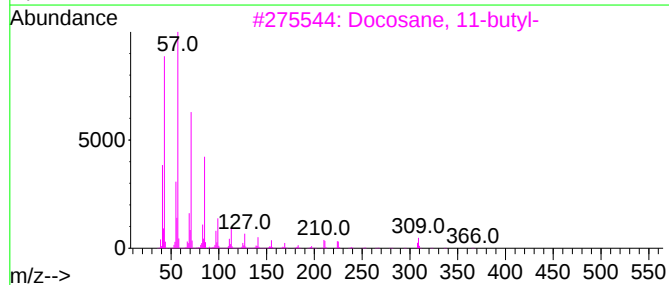
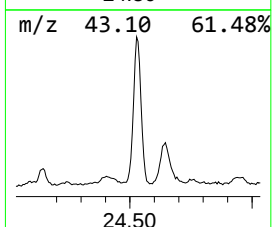
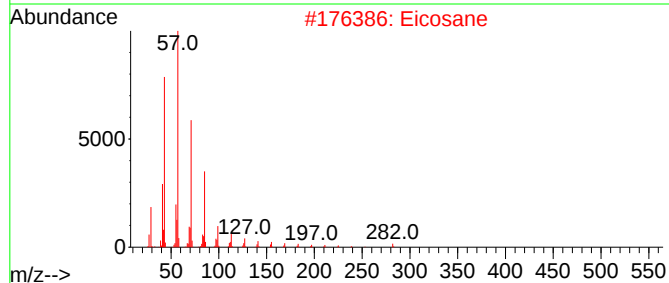
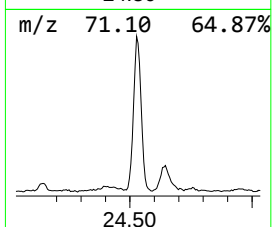
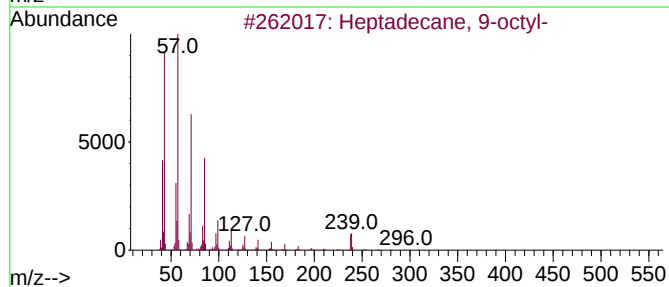
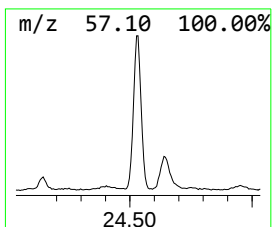
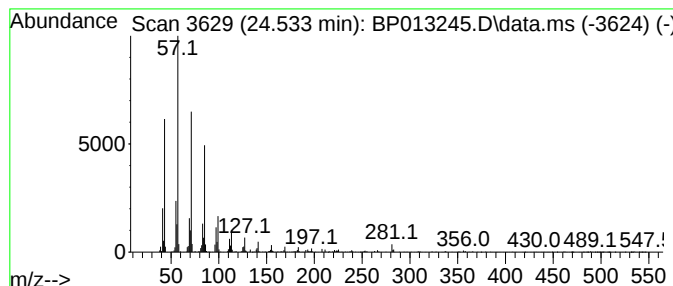
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 (DEL) Alkane: Straight-Chai... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.	
24.533	6.36 ng/ul	2000180	Perylene-d12	23.910	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptadecane, 9-octyl-	352	C25H52	007225-64-1	94
2	Eicosane	282	C20H42	000112-95-8	94
3	Docosane, 11-butyl-	366	C26H54	013475-76-8	93
4	Heneicosane, 11-pentyl-	366	C26H54	014739-72-1	93
5	Eicosane	282	C20H42	000112-95-8	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122222\
Data File : BP013245.D
Acq On : 22 Dec 2022 14:34
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Sample : N6015-04
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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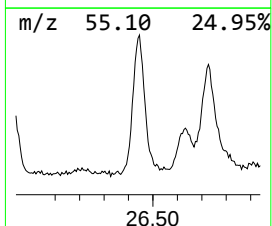
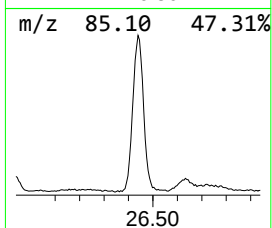
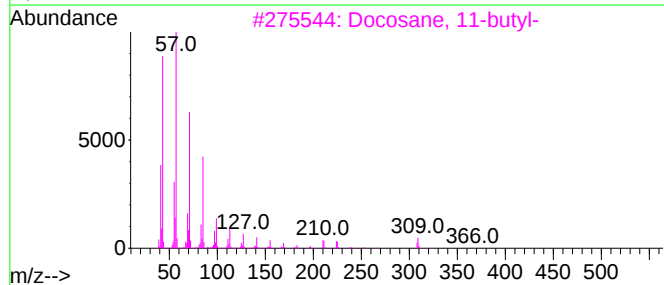
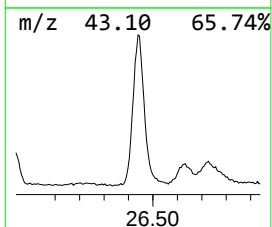
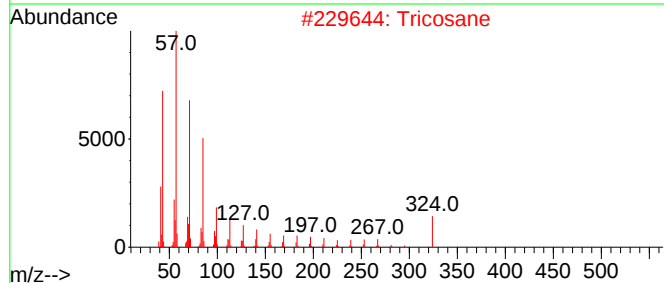
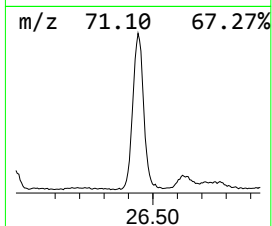
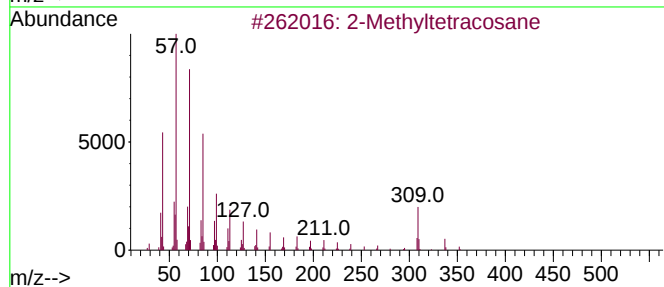
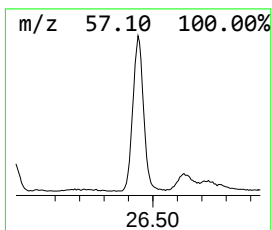
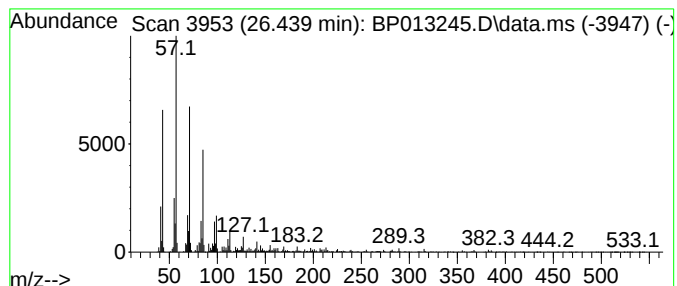
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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
26.439	14.63 ng/ul	4603040	Perylene-d12	23.910

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		2-Methyltetracosane	352	C25H52	001560-78-7	95
2		Tricosane	324	C23H48	000638-67-5	94
3		Docosane, 11-butyl-	366	C26H54	013475-76-8	93
4		Triacontane	422	C30H62	000638-68-6	93
5		Octadecane, 2-methyl-	268	C19H40	001560-88-9	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122222\
Data File : BP013245.D
Acq On : 22 Dec 2022 14:34
Operator : CG/JU
Sample : N6015-04
Misc :
ALS Vial : 12 Sample Multiplier: 1

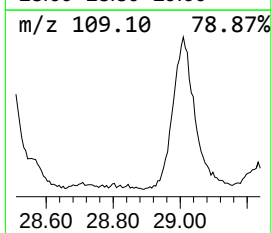
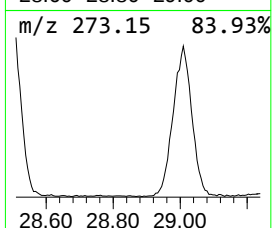
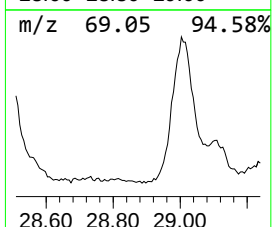
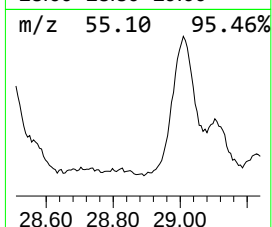
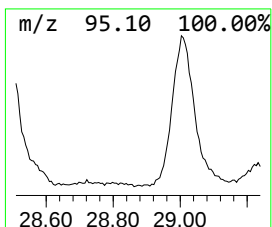
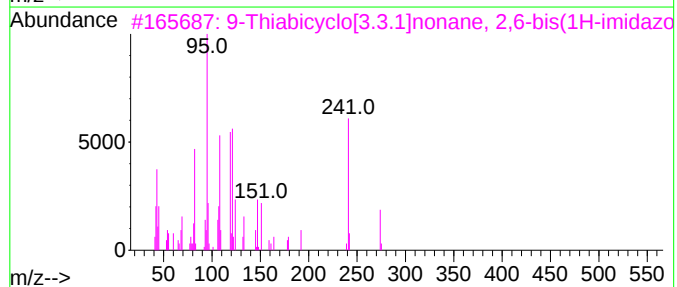
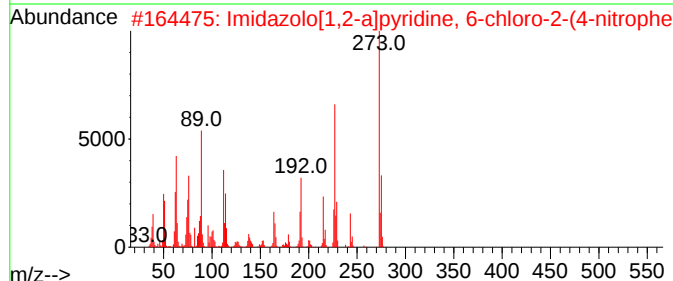
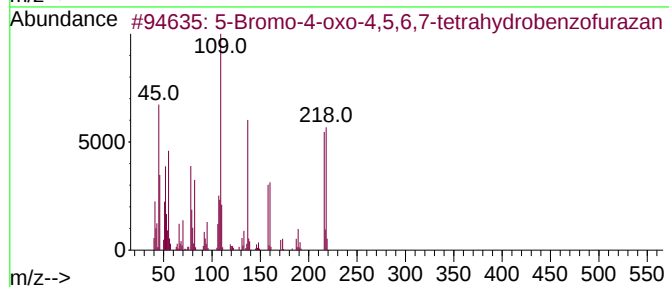
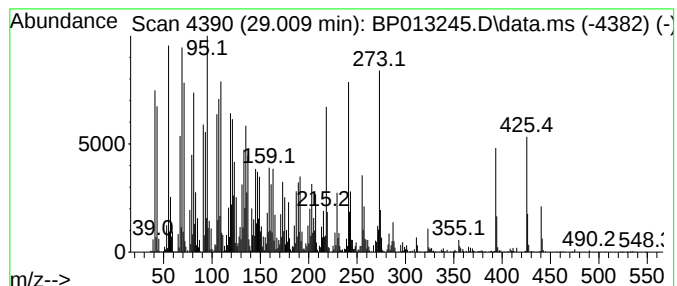
Instrument :
BNA_P
ClientSampleId :
DBXU0

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 5-Bromo-4-oxo-4,5,6,7-tetra... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.	
29.009	17.41 ng/ul	5476790	Perylene-d12	23.910	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	5-Bromo-4-oxo-4,5,6,7-tetrahydro...	216	C6H5BrN2O2	300574-36-1	91
2	Imidazolo[1,2-a]pyridine, 6-chlo...	273	C13H8ClN3O2	118000-62-7	56
3	9-Thiabicyclo[3.3.1]nonane, 2,6-...	274	C14H18N4S	1000141-52-1	38
4	Pregnan-18-ol, (5.alpha.)-	304	C21H36O	056053-09-9	30
5	Lupeol	426	C30H50O	000545-47-1	30



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122222\
Data File : BP013245.D
Acq On : 22 Dec 2022 14:34
Operator : CG/JU
Sample : N6015-04
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DBXU0

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	15.951	11.0	ng/ul	3837400	3	14.593	6998700	20.0
n-Hexadecanoic ...	18.222	5.5	ng/ul	2306650	4	17.351	8461990	20.0
Benzene, hexaet...	18.787	3.8	ng/ul	1617540	4	17.351	8461990	20.0
Octadecanoic acid	19.451	2.8	ng/ul	832040	5	21.433	5957300	20.0
4-(3-Fluoro-8-n...	19.828	35.1	ng/ul	10450600	5	21.433	5957300	20.0
3,5-di-tert-But...	20.175	6.1	ng/ul	1818730	5	21.433	5957300	20.0
Heptafluorobuty...	21.128	7.7	ng/ul	2299910	5	21.433	5957300	20.0
(DEL) Alkane: S...	23.127	2.8	ng/ul	893865	6	23.910	6293270	20.0
(DEL) Alkane: S...	24.533	6.4	ng/ul	2000180	6	23.910	6293270	20.0
(DEL) Alkane: S...	26.439	14.6	ng/ul	4603040	6	23.910	6293270	20.0
5-Bromo-4-oxo-4...	29.009	17.4	ng/ul	5476790	6	23.910	6293270	20.0