

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062323\
 Data File : BP015676.D
 Acq On : 23 Jun 2023 17:46
 Operator : MA/JU
 Sample : 03084-14
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 DCFN1

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-BP060723.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BP015676.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.399	32	36	46	rVB	30562	44687	1.29%	0.100%
2	3.846	108	112	115	rBV	18619	23059	0.67%	0.052%
3	3.875	115	117	123	rVV2	16305	27447	0.79%	0.062%
4	3.946	125	129	136	rVV	122760	162520	4.69%	0.365%
5	4.022	139	142	146	rVV2	13434	18575	0.54%	0.042%
6	4.070	146	150	157	rVB	23987	33910	0.98%	0.076%
7	4.170	163	167	172	rVB	17232	24175	0.70%	0.054%
8	4.752	261	266	273	rVB	30740	43391	1.25%	0.097%
9	5.234	340	348	354	rBV4	14171	28950	0.84%	0.065%
10	6.046	480	486	493	rBV6	12421	27489	0.79%	0.062%
11	7.246	679	690	710	rVV	327334	684363	19.77%	1.536%
12	7.405	710	717	729	rVV	290086	499158	14.42%	1.121%
13	7.605	745	751	760	rBV	395616	689324	19.91%	1.547%
14	7.675	760	763	767	rVV	18451	28692	0.83%	0.064%
15	7.716	767	770	778	rVB3	10361	17681	0.51%	0.040%
16	8.075	825	831	846	rVB	639608	1096863	31.68%	2.462%
17	8.628	919	925	934	rVB3	14729	32770	0.95%	0.074%
18	8.805	946	955	969	rBV	314320	629011	18.17%	1.412%
19	8.922	969	975	984	rVB2	97236	175913	5.08%	0.395%
20	9.263	1025	1033	1048	rBV	367424	705047	20.36%	1.583%
21	9.416	1057	1059	1070	rVB	7376	17479	0.50%	0.039%
22	9.993	1149	1157	1167	rBV	310540	593472	17.14%	1.332%
23	10.204	1185	1193	1199	rBV4	15842	42967	1.24%	0.096%
24	10.357	1215	1219	1231	rVB6	13087	30092	0.87%	0.068%
25	10.546	1243	1251	1273	rBV	318530	789804	22.81%	1.773%
26	10.904	1305	1312	1317	rVV	866209	1499048	43.30%	3.365%
27	10.957	1317	1321	1331	rVV	227832	397128	11.47%	0.891%
28	11.075	1334	1341	1358	rVB2	53964	156173	4.51%	0.351%
29	12.298	1545	1549	1560	rBV4	14823	34414	0.99%	0.077%
30	12.569	1588	1595	1610	rVB	63813	128092	3.70%	0.288%
31	12.781	1627	1631	1639	rVB	27935	47524	1.37%	0.107%
32	13.240	1705	1709	1715	rVV	13816	18004	0.52%	0.040%
33	13.528	1753	1758	1761	rBV2	19186	29076	0.84%	0.065%
34	13.569	1761	1765	1768	rVV2	28625	49678	1.43%	0.112%
35	13.604	1768	1771	1779	rVV4	26535	44049	1.27%	0.099%
36	13.887	1815	1819	1830	rBV6	18075	55456	1.60%	0.124%

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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

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Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP060723.MA.M
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37	14.051	1841	1847	1852	rBV3	20884	39387	1.14%	0.088%
38	14.134	1852	1861	1867	rBV	787749	1139235	32.90%	2.557%
39	14.181	1867	1869	1875	rVB4	26145	38450	1.11%	0.086%
40	14.287	1882	1887	1896	rVB4	10397	26805	0.77%	0.060%
41	14.422	1903	1910	1914	rBV	843142	1369119	39.54%	3.073%
42	14.598	1936	1940	1948	rVB2	23539	34285	0.99%	0.077%
43	14.728	1950	1962	1970	rBV	1171938	1763898	50.95%	3.960%
44	14.792	1970	1973	1980	rVB2	34037	60538	1.75%	0.136%
45	14.981	2000	2005	2022	rVV2	78073	271875	7.85%	0.610%
46	15.134	2025	2031	2040	rVV	90892	175920	5.08%	0.395%
47	15.204	2040	2043	2052	rVB9	15027	34346	0.99%	0.077%
48	15.340	2061	2066	2073	rBV9	12052	27380	0.79%	0.061%
49	15.510	2090	2095	2099	rBV7	13759	27953	0.81%	0.063%
50	15.551	2099	2102	2107	rVB	28989	36832	1.06%	0.083%
51	15.728	2124	2132	2138	rBV	911117	1437830	41.53%	3.228%
52	15.781	2138	2141	2151	rVB2	42143	76504	2.21%	0.172%
53	15.875	2151	2157	2165	rBV	168892	296464	8.56%	0.666%
54	16.087	2187	2193	2206	rVB2	113627	211055	6.10%	0.474%
55	16.228	2212	2217	2224	rBV2	19662	37500	1.08%	0.084%
56	16.428	2245	2251	2255	rBV3	61221	143744	4.15%	0.323%
57	16.469	2255	2258	2265	rVB2	54090	91098	2.63%	0.204%
58	16.657	2287	2290	2297	rVB4	16551	25093	0.72%	0.056%
59	17.016	2348	2351	2356	rBV7	16376	22897	0.66%	0.051%
60	17.151	2369	2374	2380	rBV3	61403	106116	3.06%	0.238%
61	17.216	2381	2385	2390	rBV	47810	67781	1.96%	0.152%
62	17.310	2397	2401	2407	rVB	29823	44958	1.30%	0.101%
63	17.363	2407	2410	2417	rBV6	25026	44983	1.30%	0.101%
64	17.492	2426	2432	2436	rBV	1091879	1620396	46.80%	3.638%
65	17.534	2436	2439	2444	rVV	489240	670879	19.38%	1.506%
66	17.592	2444	2449	2453	rVV2	811925	1189546	34.36%	2.670%
67	17.628	2453	2455	2460	rVB	168741	217104	6.27%	0.487%
68	17.904	2500	2502	2507	rVV3	25887	32996	0.95%	0.074%
69	17.951	2507	2510	2516	rVB2	38240	49296	1.42%	0.111%
70	18.292	2564	2568	2573	rBV3	42758	67702	1.96%	0.152%
71	18.345	2573	2577	2583	rBV2	484009	693369	20.03%	1.556%
72	18.398	2583	2586	2592	rVB2	92055	158428	4.58%	0.356%
73	18.480	2597	2600	2605	rVB	41535	52357	1.51%	0.118%
74	18.539	2605	2610	2614	rBV2	108725	199567	5.76%	0.448%
75	18.622	2621	2624	2633	rBV4	40095	69505	2.01%	0.156%
76	18.833	2657	2660	2664	rVB	52575	62677	1.81%	0.141%

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77	18.886	2666	2669	2672	rBV4	56360	82483	2.38%	0.185%
78	19.228	2725	2727	2732	rVB2	125510	151947	4.39%	0.341%
79	19.492	2767	2772	2778	rVB	1467904	1983072	57.28%	4.452%
80	19.575	2783	2786	2793	rVV3	138129	206199	5.96%	0.463%
81	19.639	2794	2797	2804	rVB	137754	175152	5.06%	0.393%
82	19.822	2823	2828	2830	rBV2	1057265	1558171	45.00%	3.498%
83	19.845	2830	2832	2840	rVB	1208569	1502111	43.39%	3.372%
84	19.916	2841	2844	2848	rBV	69067	94619	2.73%	0.212%
85	20.304	2907	2910	2912	rBV3	104808	134477	3.88%	0.302%
86	20.427	2928	2931	2934	rBV	115427	126696	3.66%	0.284%
87	20.622	2961	2964	2968	rBV	117578	167990	4.85%	0.377%
88	21.063	3036	3039	3046	rBV	196960	288847	8.34%	0.648%
89	21.222	3063	3066	3068	rBV2	218823	284608	8.22%	0.639%
90	21.257	3069	3072	3076	rVB2	348637	470668	13.59%	1.057%
91	21.357	3087	3089	3098	rVB	216749	275788	7.97%	0.619%
92	21.580	3120	3127	3131	rBV2	1613959	3462262	100.00%	7.772%
93	21.622	3131	3134	3143	rVB	921832	1189844	34.37%	2.671%
94	23.292	3415	3418	3423	rVB	302125	422525	12.20%	0.948%
95	23.363	3423	3430	3435	rBV	1118405	2892147	83.53%	6.492%
96	23.916	3519	3524	3529	rBV	587792	1150954	33.24%	2.584%
97	23.974	3529	3534	3539	rVV2	814924	1560540	45.07%	3.503%
98	24.027	3539	3543	3554	rVB	871557	1722574	49.75%	3.867%
99	24.139	3557	3562	3568	rBV	815217	1620535	46.81%	3.638%
100	24.745	3661	3665	3679	rVB	604792	1361289	39.32%	3.056%

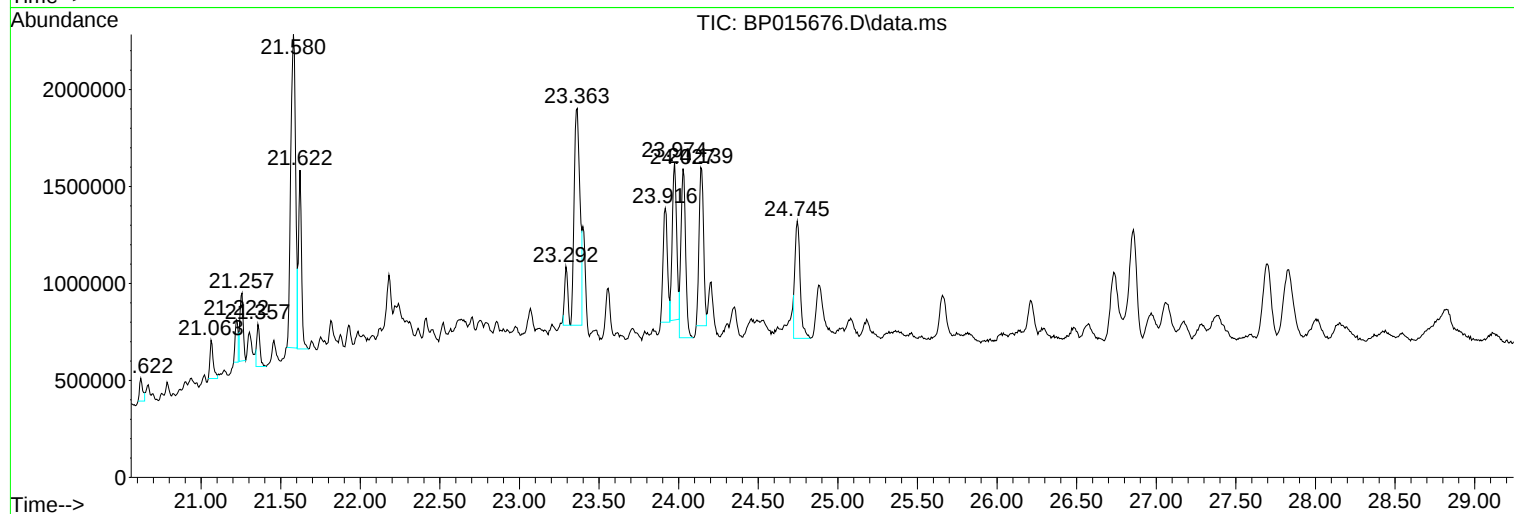
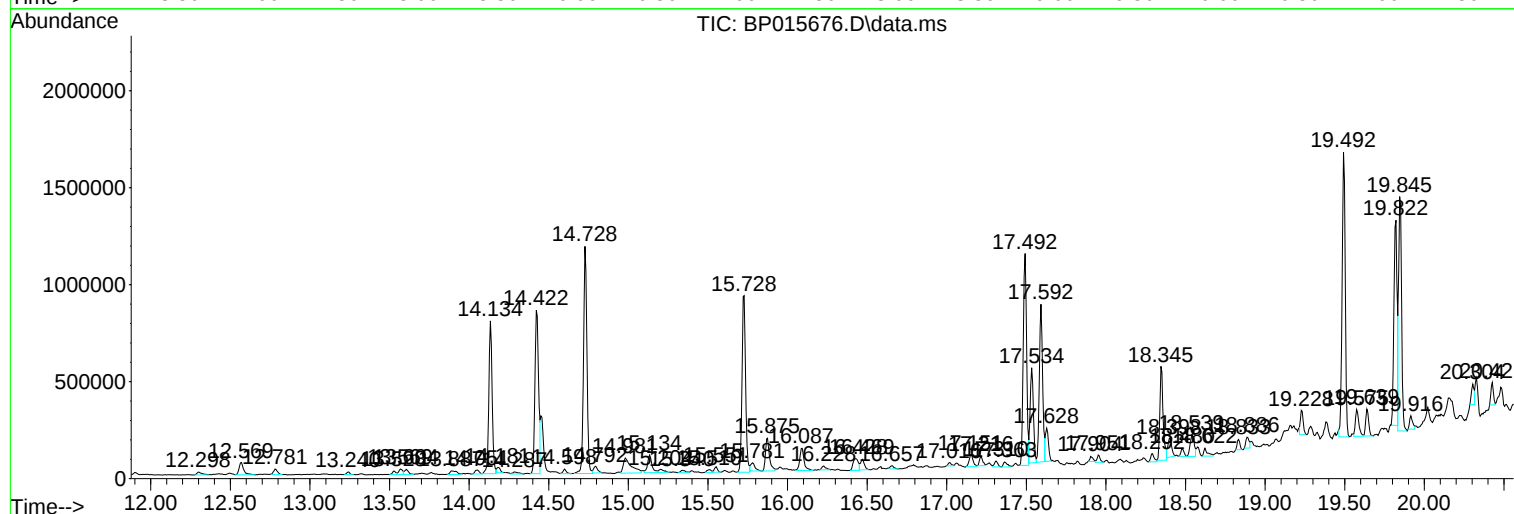
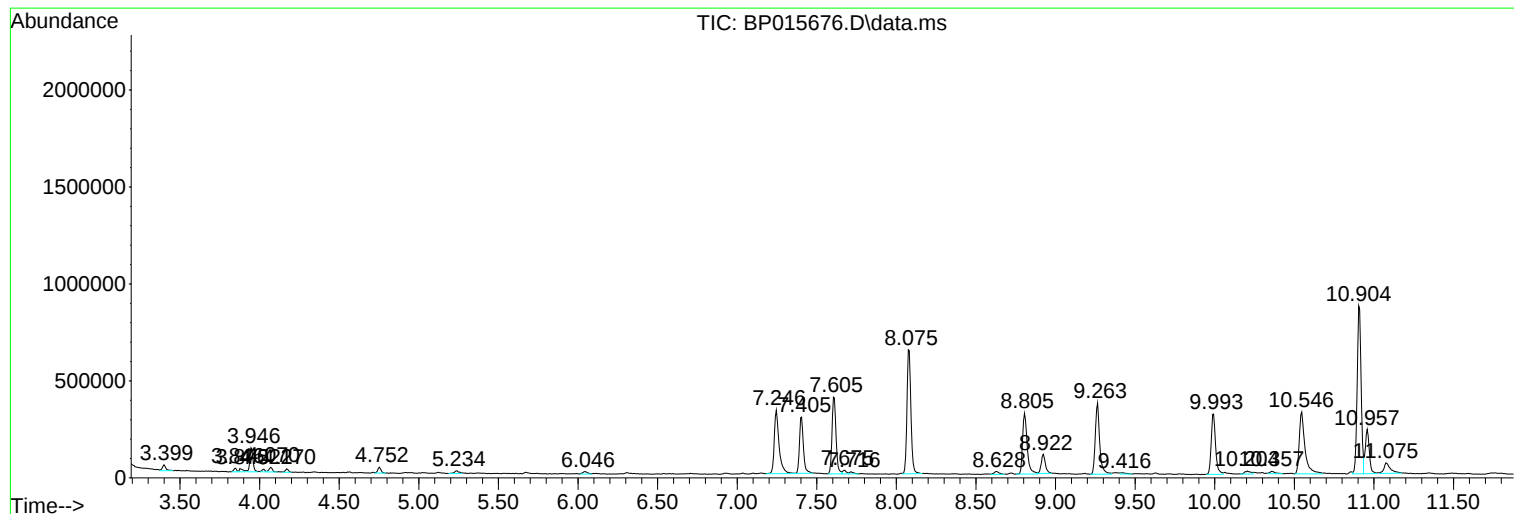
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Instrument :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP060723.MA.M
Quant Title : SVOA CALIBRATION

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



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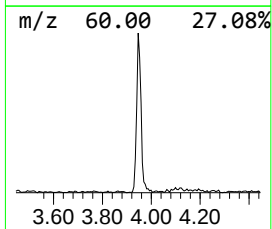
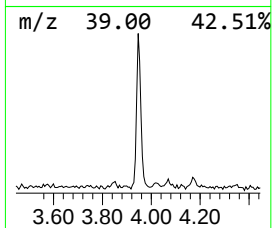
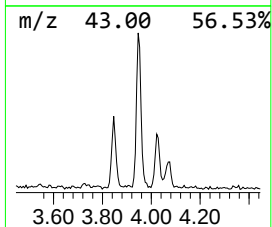
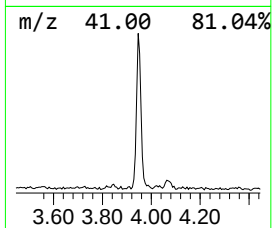
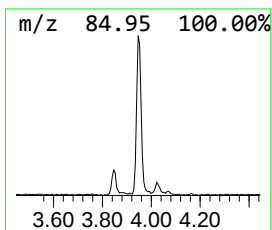
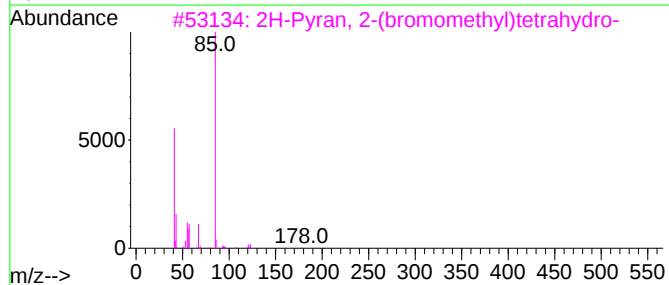
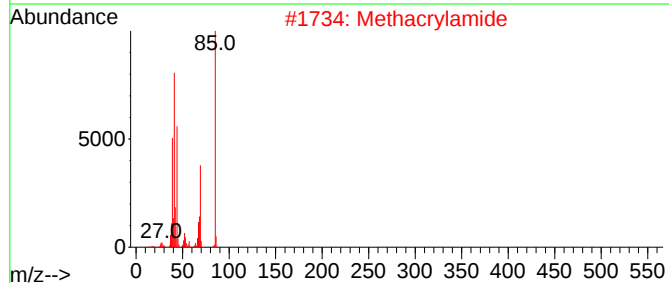
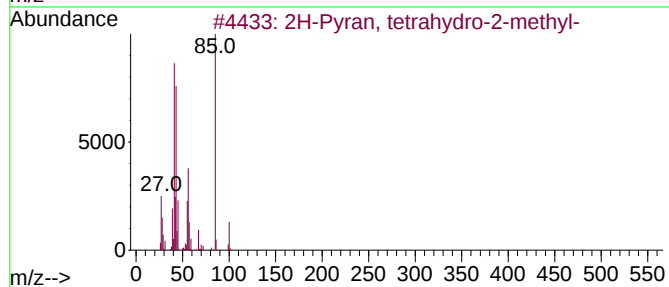
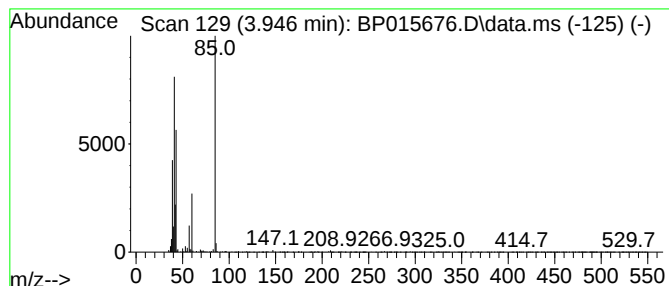
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP060723.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 2H-Pyran, tetrahydro-2-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.946	2.96 ng/ul	162520	1,4-Dichlorobenzene-d4	8.081

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2H-Pyran, tetrahydro-2-methyl-			100	C6H12O	010141-72-7	50
2	Methacrylamide			85	C4H7NO	000079-39-0	50
3	2H-Pyran, 2-(bromomethyl)tetrahy...			178	C6H11BrO	034723-82-5	40
4	Butanoyl chloride, 3-methyl-			120	C5H9ClO	000108-12-3	38
5	2H-Pyran, tetrahydro-2-methyl-			100	C6H12O	010141-72-7	33



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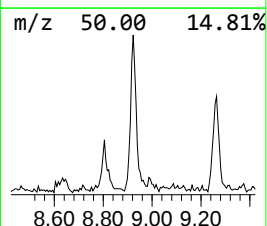
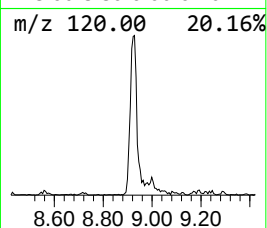
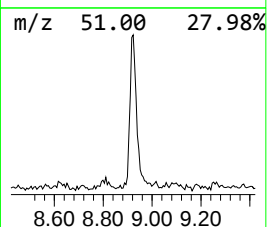
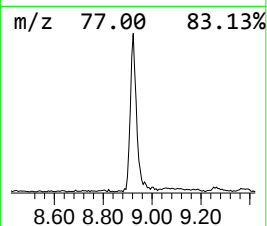
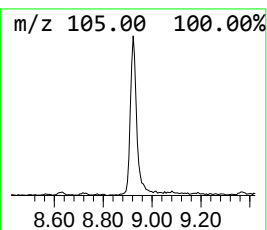
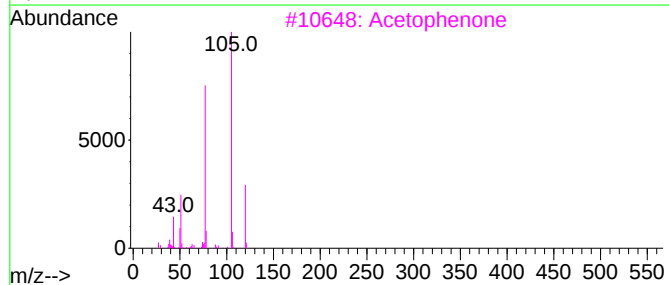
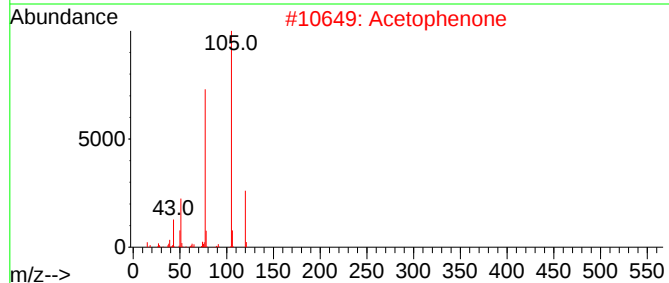
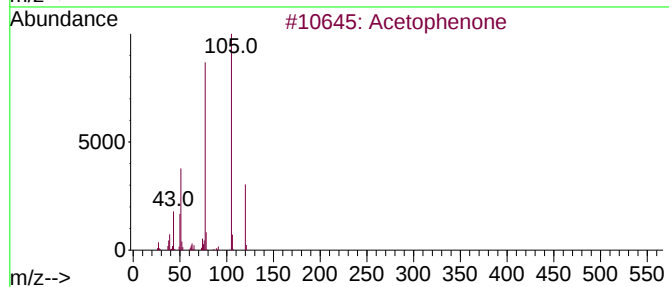
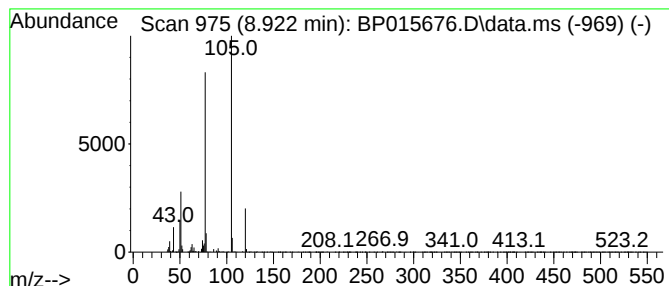
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP060723.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 Aceto phenone Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.922	3.21 ng/ul	175913	1,4-Dichlorobenzene-d4	8.081

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Acetophenone	120	C8H8O	000098-86-2	94
2			Acetophenone	120	C8H8O	000098-86-2	91
3			Acetophenone	120	C8H8O	000098-86-2	91
4			Benzoyl isothiocyanate	163	C8H5NOS	000532-55-8	72
5			Ethanone, 2,2,2-trifluoro-1-phenyl-	174	C8H5F3O	000434-45-7	72



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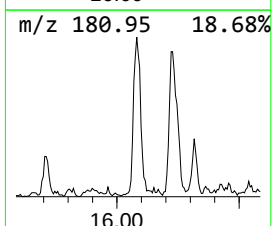
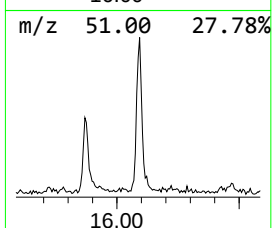
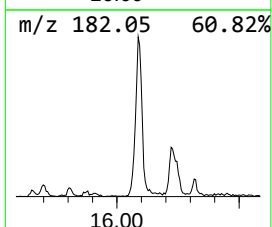
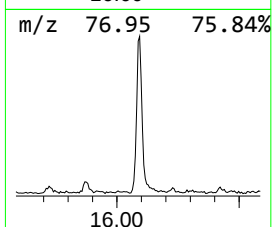
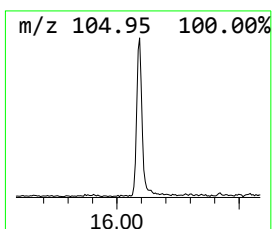
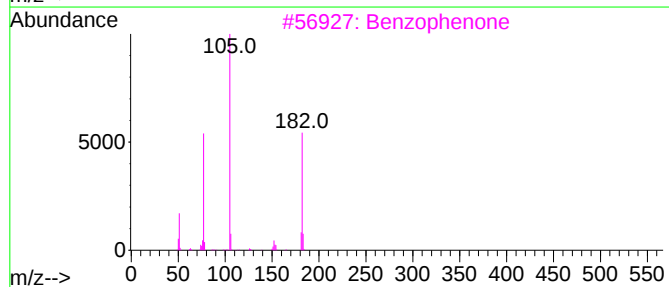
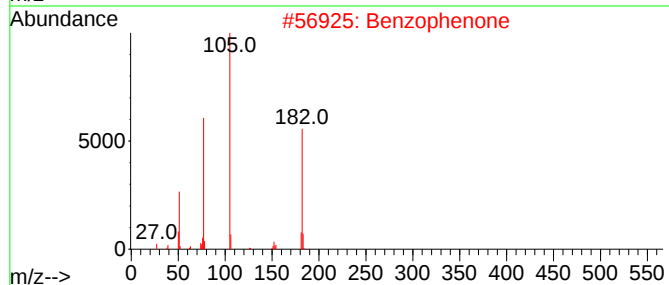
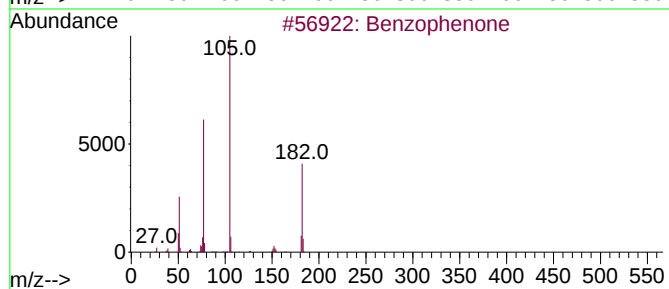
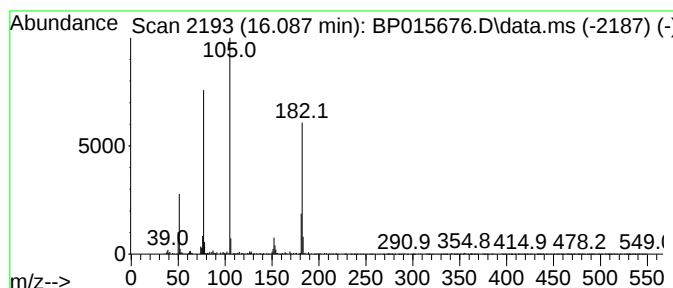
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Quant Title : SVOA CALIBRATION

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

Peak Number 3 Benzophenone Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.087	2.39 ng/ul	211055	Acenaphthene-d10	14.728

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzophenone	182	C13H10O	000119-61-9	94
2			Benzophenone	182	C13H10O	000119-61-9	94
3			Benzophenone	182	C13H10O	000119-61-9	94
4			Benzophenone	182	C13H10O	000119-61-9	86
5			Benzophenone	182	C13H10O	000119-61-9	68



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062323\
Data File : BP015676.D
Acq On : 23 Jun 2023 17:46
Operator : MA/JU
Sample : 03084-14
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCFN1

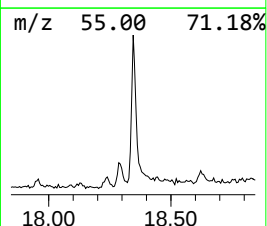
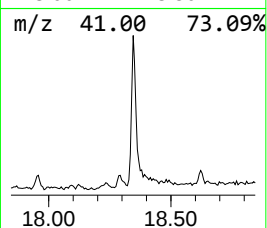
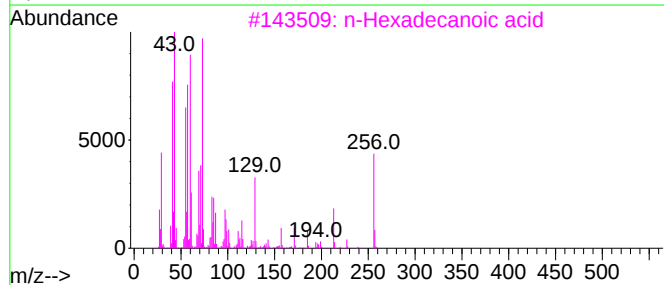
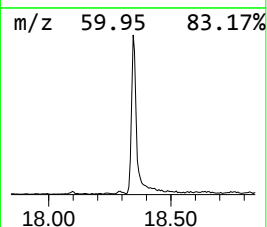
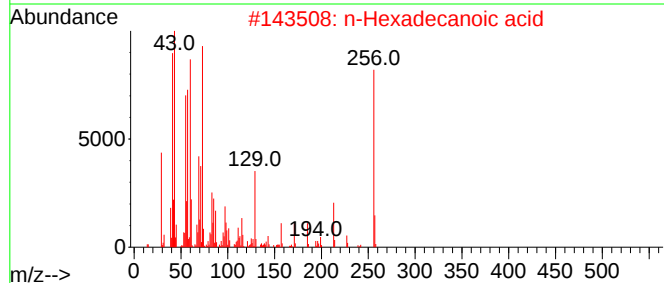
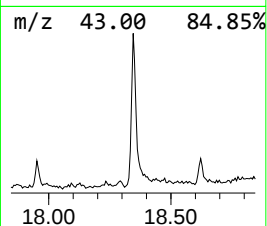
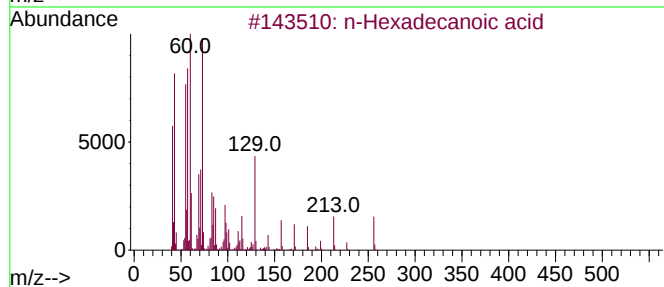
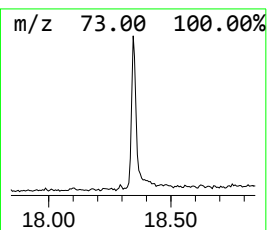
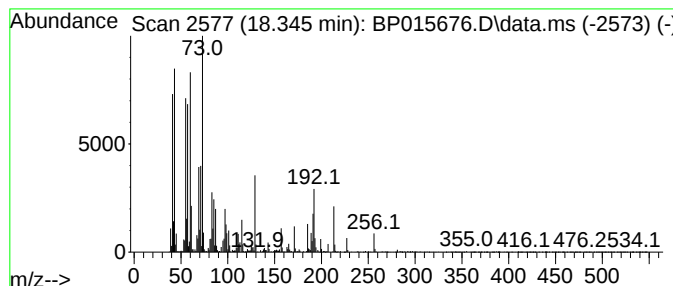
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 n-Hexadecanoic acid Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.345	8.56 ng/ul	693369	Phenanthrene-d10	17.492

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
3			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
4			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	95
5			Octadecanoic acid	284	C18H36O2	000057-11-4	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062323\
Data File : BP015676.D
Acq On : 23 Jun 2023 17:46
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Sample : 03084-14
Misc :
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Instrument :
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ClientSampleId :
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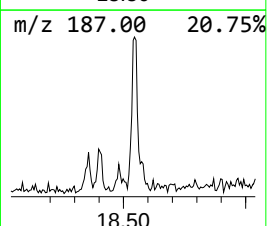
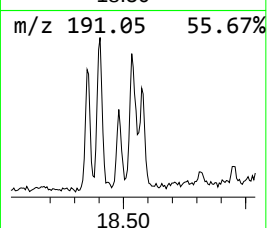
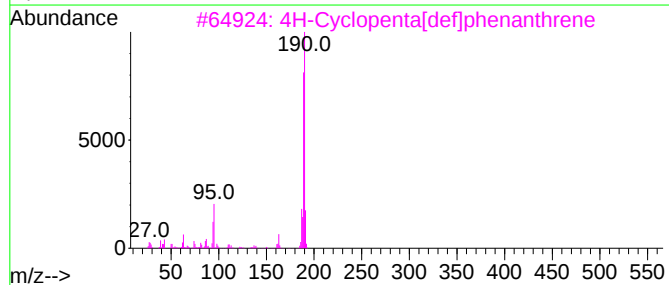
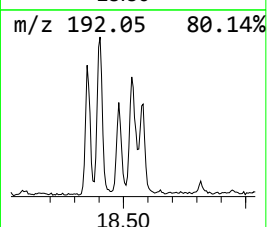
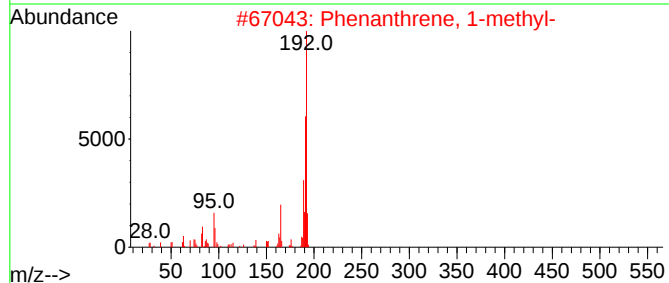
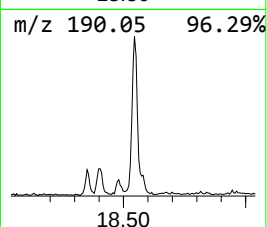
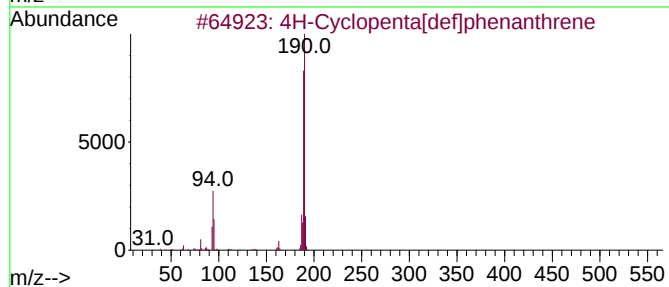
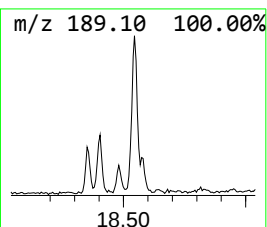
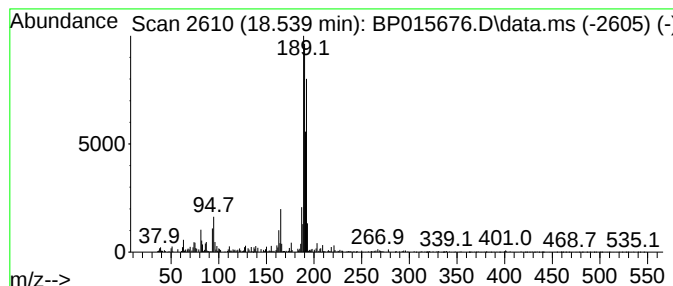
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 4H-Cyclopenta[def]phenanthrene Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.539	2.46 ng/ul	199567	Phenanthrene-d10	17.492

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	60
2			Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	53
3			4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	49
4			Carbonic acid, butyl 2-formyl-4,...	290	C12H12Cl2O4	1000331-42-2	47
5			Phenanthrene, 4-methyl-	192	C15H12	000832-64-4	46



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062323\
Data File : BP015676.D
Acq On : 23 Jun 2023 17:46
Operator : MA/JU
Sample : 03084-14
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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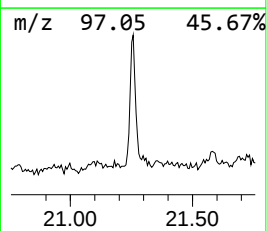
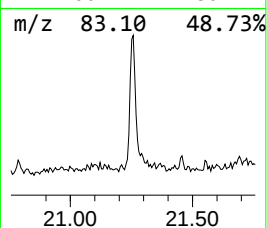
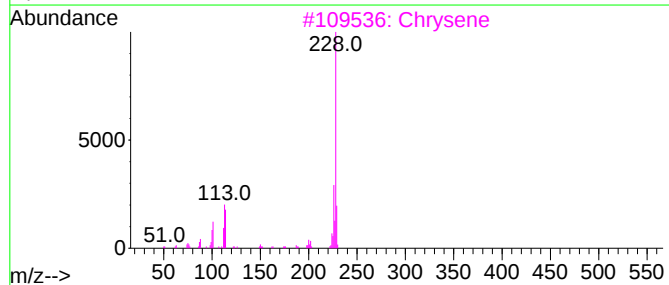
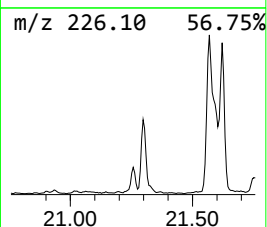
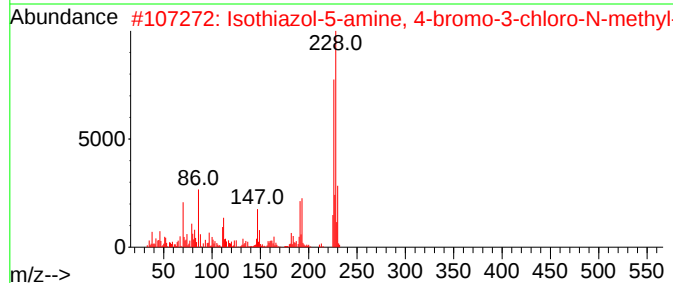
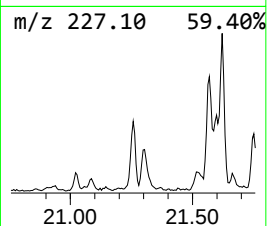
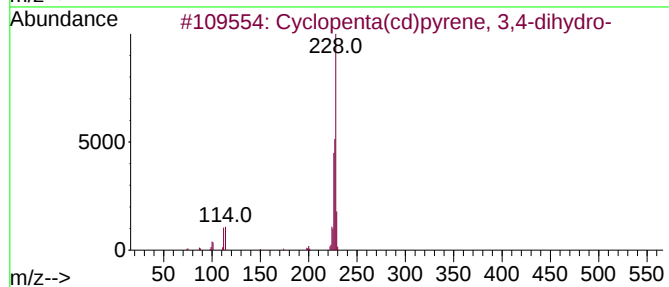
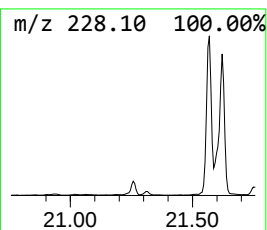
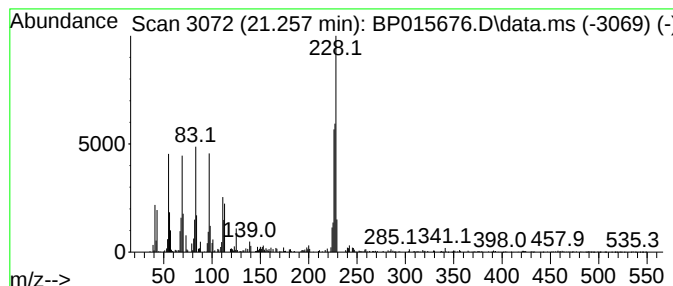
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 unknown-01 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.257	2.72 ng/ul	470668	Chrysene-d12	21.586

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclopenta(cd)pyrene, 3,4-dihydro-	228	C18H12	025732-74-5	46
2			Isothiazol-5-amine, 4-bromo-3-ch...	226	C4H4BrClN2S	1000272-59-9	43
3			Chrysene	228	C18H12	000218-01-9	42
4			Benz[a]anthracene	228	C18H12	000056-55-3	38
5			Benzo[c]phenanthrene	228	C18H12	000195-19-7	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062323\
Data File : BP015676.D
Acq On : 23 Jun 2023 17:46
Operator : MA/JU
Sample : 03084-14
Misc :
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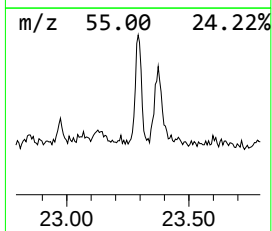
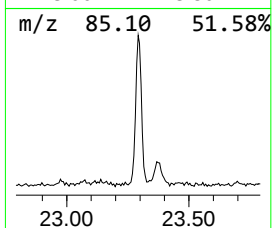
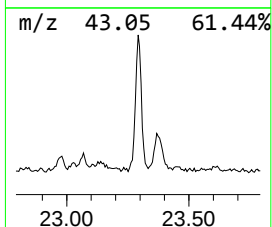
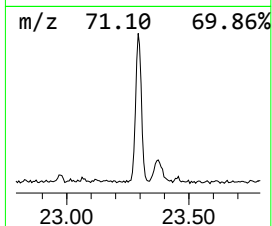
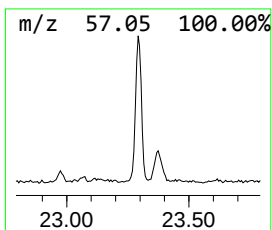
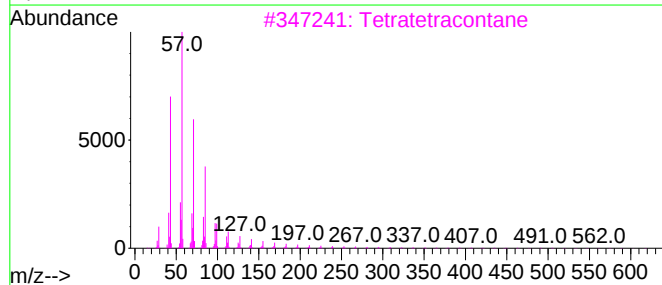
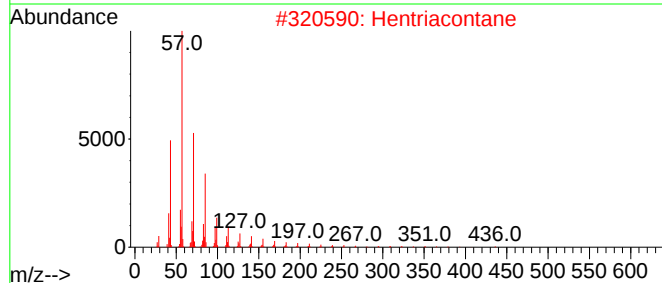
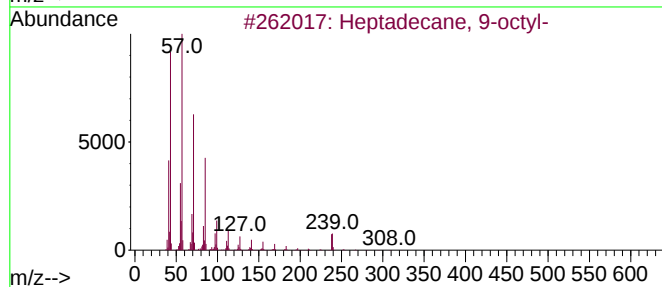
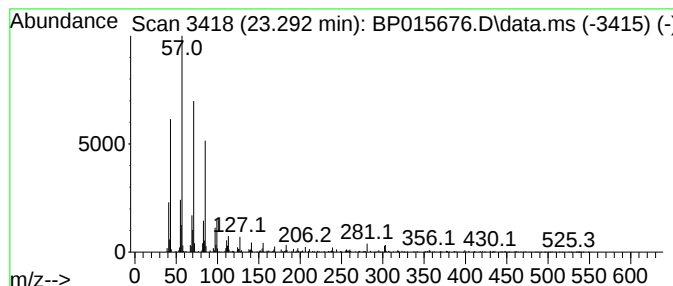
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ClientSampleId :
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 (DEL) Alkane: Straight-Chai... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.	
23.292	5.21 ng/ul	422525	Perylene-d12	24.139	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptadecane, 9-octyl-	352	C25H52	007225-64-1	93
2	Hentriacontane	437	C31H64	000630-04-6	93
3	Tetratetracontane	619	C44H90	007098-22-8	87
4	Dodecane, 2-methyl-	184	C13H28	001560-97-0	87
5	Tetracosane	338	C24H50	000646-31-1	83



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062323\
Data File : BP015676.D
Acq On : 23 Jun 2023 17:46
Operator : MA/JU
Sample : 03084-14
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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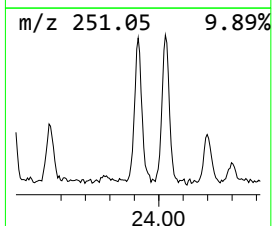
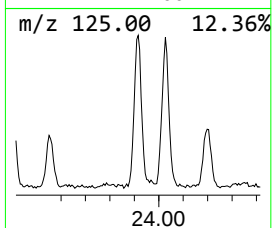
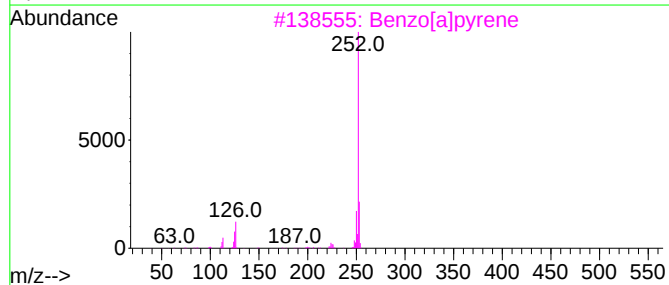
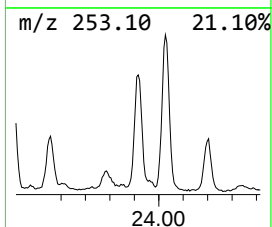
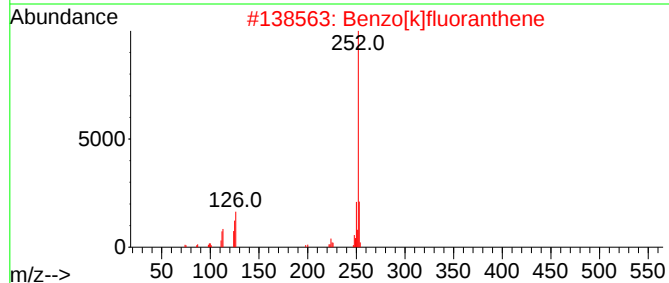
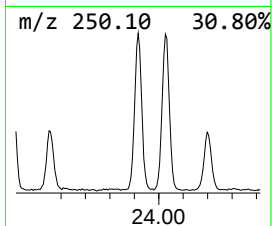
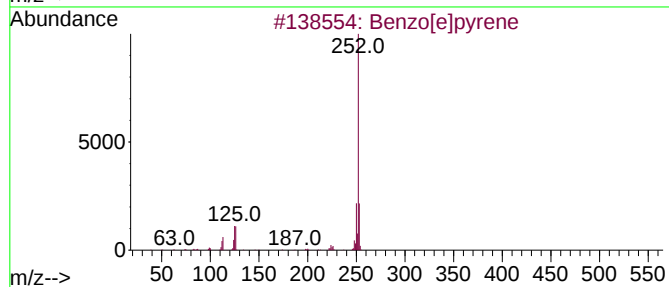
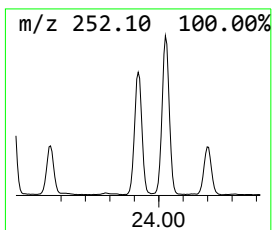
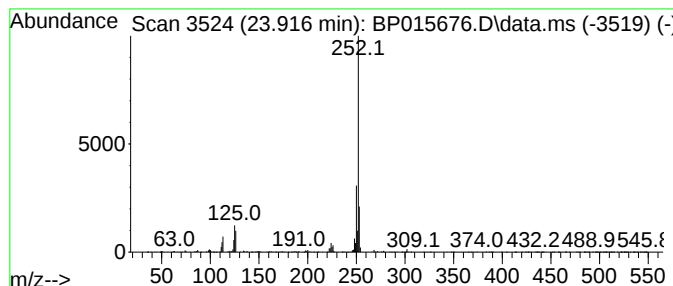
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 Benzo[e]pyrene Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.916	14.20 ng/ul	1150950	Perylene-d12	24.139

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzo[e]pyrene	252	C20H12	000192-97-2	99
2			Benzo[k]fluoranthene	252	C20H12	000207-08-9	97
3			Benzo[a]pyrene	252	C20H12	000050-32-8	95
4			Perylene	252	C20H12	000198-55-0	95
5			Benzo[b]fluoranthene	252	C20H12	000205-99-2	95



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062323\
Data File : BP015676.D
Acq On : 23 Jun 2023 17:46
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Sample : 03084-14
Misc :
ALS Vial : 10 Sample Multiplier: 1

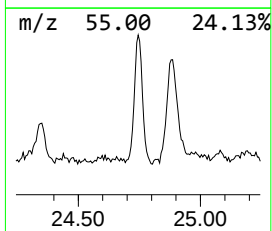
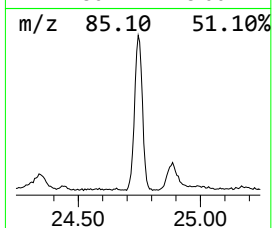
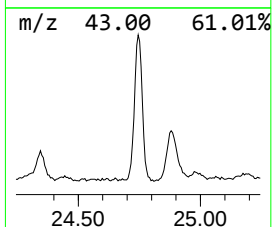
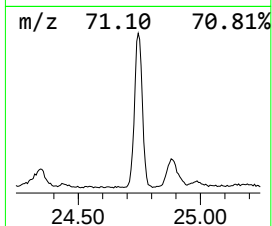
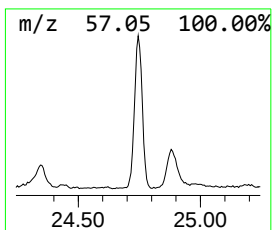
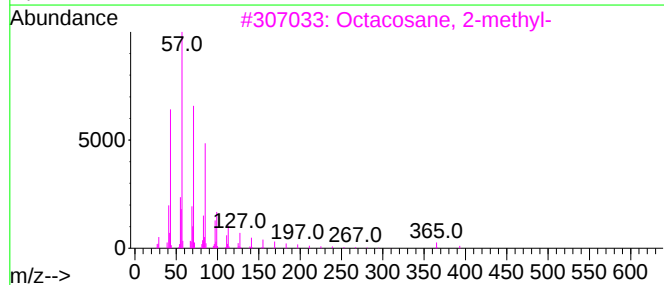
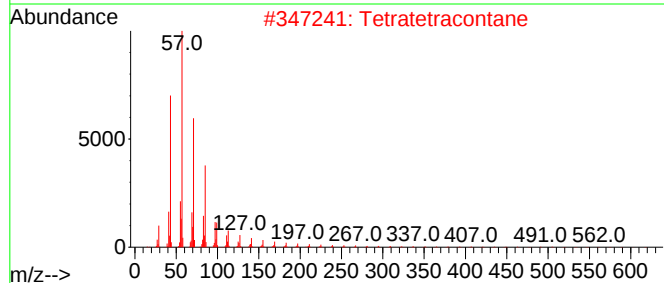
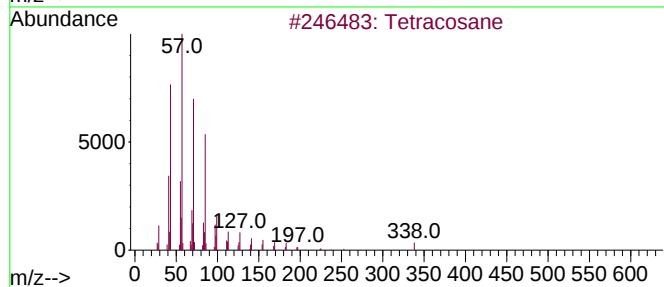
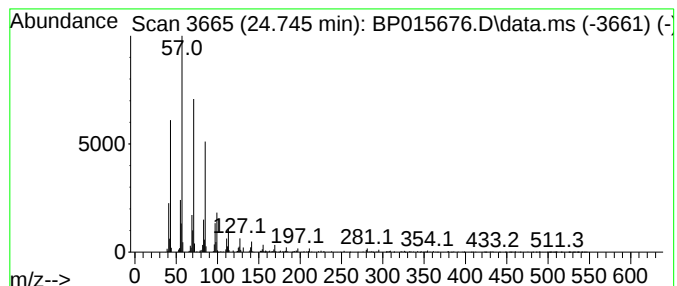
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP060723.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 9 (DEL) Alkane: Straight-Chai... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.	
24.745	16.80 ng/ul	1361290	Perylene-d12	24.139	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tetracosane	338	C24H50	000646-31-1	96
2	Tetratetracontane	619	C44H90	007098-22-8	91
3	Octacosane, 2-methyl-	408	C29H60	001560-98-1	90
4	Eicosane, 7-hexyl-	366	C26H54	055333-99-8	87
5	Dotriacontane, 1-iodo-	576	C32H65I	1000406-32-4	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062323\
Data File : BP015676.D
Acq On : 23 Jun 2023 17:46
Operator : MA/JU
Sample : 03084-14
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCFN1

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP060723.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST0.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
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Aceto phenone	8.922	3.2	ng/ul	175913	1	8.081	1096860	20.0
Benzophenone	16.087	2.4	ng/ul	211055	3	14.728	1763900	20.0
n-Hexadecanoic ...	18.345	8.6	ng/ul	693369	4	17.492	1620400	20.0
4H-Cyclopenta[d...	18.539	2.5	ng/ul	199567	4	17.492	1620400	20.0
unknown-01	21.257	2.7	ng/ul	470668	5	21.586	3462260	20.0
(DEL) Alkane: S...	23.292	5.2	ng/ul	422525	6	24.139	1620540	20.0
Benzo[e]pyrene	23.916	14.2	ng/ul	1150950	6	24.139	1620540	20.0
(DEL) Alkane: S...	24.745	16.8	ng/ul	1361290	6	24.139	1620540	20.0