

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP050423\
 Data File : BP014967.D
 Acq On : 04 May 2023 23:21
 Operator : CG/JU
 Sample : 02479-06
 Misc :
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 EW947

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-BP050323.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BP014967.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.464	10	13	24	rVB2	38788	82724	1.77%	0.167%
2	3.911	86	89	93	rVB	36543	45354	0.97%	0.091%
3	4.017	102	107	115	rVB	275254	350551	7.50%	0.707%
4	4.093	116	120	124	rVV	29214	37345	0.80%	0.075%
5	4.146	125	129	134	rVB	31762	46919	1.00%	0.095%
6	4.246	142	146	151	rVB	28189	37918	0.81%	0.076%
7	4.634	207	212	216	rVB5	14360	23267	0.50%	0.047%
8	4.828	241	245	255	rVB	40912	61879	1.32%	0.125%
9	5.199	303	308	314	rBV	18146	26328	0.56%	0.053%
10	5.299	319	325	331	rBV2	52913	81654	1.75%	0.165%
11	5.822	409	414	418	rVB	19776	29097	0.62%	0.059%
12	6.140	460	468	472	rBV5	18020	31235	0.67%	0.063%
13	7.128	627	636	643	rBV3	19974	36070	0.77%	0.073%
14	7.311	659	667	675	rBV	422240	724730	15.50%	1.461%
15	7.481	690	696	703	rBV	403846	617009	13.19%	1.244%
16	7.687	725	731	741	rVB	444407	721614	15.43%	1.455%
17	7.775	741	746	748	rBV	20984	28774	0.62%	0.058%
18	7.805	748	751	755	rVB	25170	34296	0.73%	0.069%
19	8.169	803	813	820	rBV	1075279	1738256	37.17%	3.505%
20	8.281	828	832	839	rVB4	12450	22467	0.48%	0.045%
21	8.716	903	906	917	rVB5	19958	44366	0.95%	0.089%
22	8.811	917	922	927	rBV4	24531	47281	1.01%	0.095%
23	8.875	927	933	941	rVV	402991	661173	14.14%	1.333%
24	8.999	948	954	961	rVB	48029	83192	1.78%	0.168%
25	9.281	993	1002	1006	rBV2	28896	58978	1.26%	0.119%
26	9.340	1006	1012	1019	rVV	384816	665073	14.22%	1.341%
27	9.399	1019	1022	1027	rVV2	24796	35774	0.77%	0.072%
28	9.534	1041	1045	1053	rVB6	10809	24414	0.52%	0.049%
29	9.740	1075	1080	1087	rBV5	10769	24930	0.53%	0.050%
30	9.810	1087	1092	1098	rVB3	16254	28203	0.60%	0.057%
31	9.875	1098	1103	1107	rBV	18048	30992	0.66%	0.062%
32	10.069	1129	1136	1148	rBV	292548	521624	11.15%	1.052%
33	10.193	1149	1157	1158	rBV6	12774	25869	0.55%	0.052%
34	10.393	1186	1191	1198	rBV4	27301	61507	1.32%	0.124%
35	10.616	1222	1229	1236	rVV	431215	734820	15.71%	1.482%
36	10.693	1236	1242	1250	rVB2	16694	34761	0.74%	0.070%

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Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP050323.MA.M
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37	11.005	1282	1295	1300	rBV	1276127	2263775	48.41%	4.564%
38	11.052	1300	1303	1310	rVV	179021	286920	6.14%	0.578%
39	11.140	1310	1318	1328	rVV	162584	346574	7.41%	0.699%
40	11.704	1409	1414	1419	rVB7	12661	24049	0.51%	0.048%
41	11.769	1420	1425	1430	rBV9	12754	23531	0.50%	0.047%
42	11.904	1442	1448	1455	rBV9	16752	35341	0.76%	0.071%
43	12.004	1461	1465	1470	rBV3	19642	31367	0.67%	0.063%
44	12.399	1527	1532	1539	rBV2	34343	52633	1.13%	0.106%
45	12.646	1565	1574	1580	rVV	381669	594951	12.72%	1.200%
46	12.699	1580	1583	1589	rVB2	30922	41288	0.88%	0.083%
47	12.863	1604	1611	1619	rVB	177886	299420	6.40%	0.604%
48	13.081	1641	1648	1653	rBV9	9077	25767	0.55%	0.052%
49	13.340	1685	1692	1698	rVB4	32653	58564	1.25%	0.118%
50	13.622	1733	1740	1748	rBV2	66043	134549	2.88%	0.271%
51	13.840	1772	1777	1787	rVB	68144	145057	3.10%	0.292%
52	13.969	1790	1799	1801	rBV	163823	264891	5.66%	0.534%
53	14.128	1820	1826	1831	rBV	200372	355411	7.60%	0.717%
54	14.210	1831	1840	1845	rVB	681265	1141703	24.41%	2.302%
55	14.275	1847	1851	1855	rVB4	47658	71814	1.54%	0.145%
56	14.322	1855	1859	1862	rBV4	18619	26019	0.56%	0.052%
57	14.369	1862	1867	1870	rBV	89526	134174	2.87%	0.271%
58	14.504	1884	1890	1903	rBV	860698	1422777	30.43%	2.869%
59	14.622	1906	1910	1917	rVB7	24645	44502	0.95%	0.090%
60	14.687	1917	1921	1925	rBV	61709	79531	1.70%	0.160%
61	14.810	1932	1942	1947	rBV	1570900	2391719	51.15%	4.822%
62	14.869	1947	1952	1959	rVB4	49704	131916	2.82%	0.266%
63	14.998	1968	1974	1984	rBV3	117607	283588	6.06%	0.572%
64	15.093	1987	1990	1995	rVV4	22455	31782	0.68%	0.064%
65	15.204	2003	2009	2014	rVB2	130767	214139	4.58%	0.432%
66	15.275	2017	2021	2025	rBV	87304	122068	2.61%	0.246%
67	15.422	2041	2046	2051	rBV2	65568	103308	2.21%	0.208%
68	15.475	2051	2055	2059	rVB	77926	103218	2.21%	0.208%
69	15.593	2069	2075	2078	rBV2	56677	101010	2.16%	0.204%
70	15.634	2078	2082	2086	rVB2	93130	135253	2.89%	0.273%
71	15.798	2105	2110	2115	rBV	1033935	1401983	29.98%	2.827%
72	15.845	2115	2118	2123	rVV3	46112	63401	1.36%	0.128%
73	15.910	2125	2129	2134	rVB5	95921	146127	3.12%	0.295%
74	16.157	2166	2171	2176	rVB2	268735	400082	8.56%	0.807%
75	16.292	2192	2194	2201	rVB2	36589	64810	1.39%	0.131%
76	16.504	2226	2230	2231	rBV	78117	98700	2.11%	0.199%

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77	16.528	2231	2234	2242	rVB2	196390	289454	6.19%	0.584%
78	16.663	2254	2257	2262	rBV4	33224	52074	1.11%	0.105%
79	17.210	2347	2350	2357	rBV2	75103	112079	2.40%	0.226%
80	17.298	2361	2365	2370	rVB4	84520	119925	2.56%	0.242%
81	17.351	2371	2374	2377	rBV2	50429	72690	1.55%	0.147%
82	17.434	2386	2388	2391	rBV3	32651	39160	0.84%	0.079%
83	17.563	2404	2410	2414	rBV	1720990	2520047	53.89%	5.081%
84	17.604	2414	2417	2422	rVB	754777	974807	20.85%	1.965%
85	17.663	2422	2427	2431	rBV	1125315	1557058	33.30%	3.139%
86	17.957	2474	2477	2481	rBV	75780	104855	2.24%	0.211%
87	18.034	2488	2490	2494	rVB4	76545	81085	1.73%	0.163%
88	18.139	2504	2508	2516	rVB3	157022	290116	6.20%	0.585%
89	18.416	2551	2555	2560	rBV	430329	626955	13.41%	1.264%
90	18.463	2560	2563	2567	rVB	212723	282471	6.04%	0.570%
91	18.604	2582	2587	2591	rBV2	244851	459001	9.82%	0.925%
92	19.557	2744	2749	2755	rVB	1767112	2297273	49.13%	4.632%
93	19.881	2799	2804	2806	rBV2	1748985	2427612	51.91%	4.894%
94	19.910	2806	2809	2816	rVB	1541927	2041495	43.66%	4.116%
95	20.763	2950	2954	2957	rBV	1058650	1123733	24.03%	2.266%
96	21.310	3045	3047	3052	rVB2	621239	742030	15.87%	1.496%
97	21.527	3079	3084	3092	rBV	2366616	3314427	70.88%	6.682%
98	21.645	3097	3104	3108	rVV2	2348009	4676339	100.00%	9.428%
99	21.680	3108	3110	3116	rVB	826552	990704	21.19%	1.997%
100	24.245	3541	3546	3553	rVB2	1434769	2846141	60.86%	5.738%

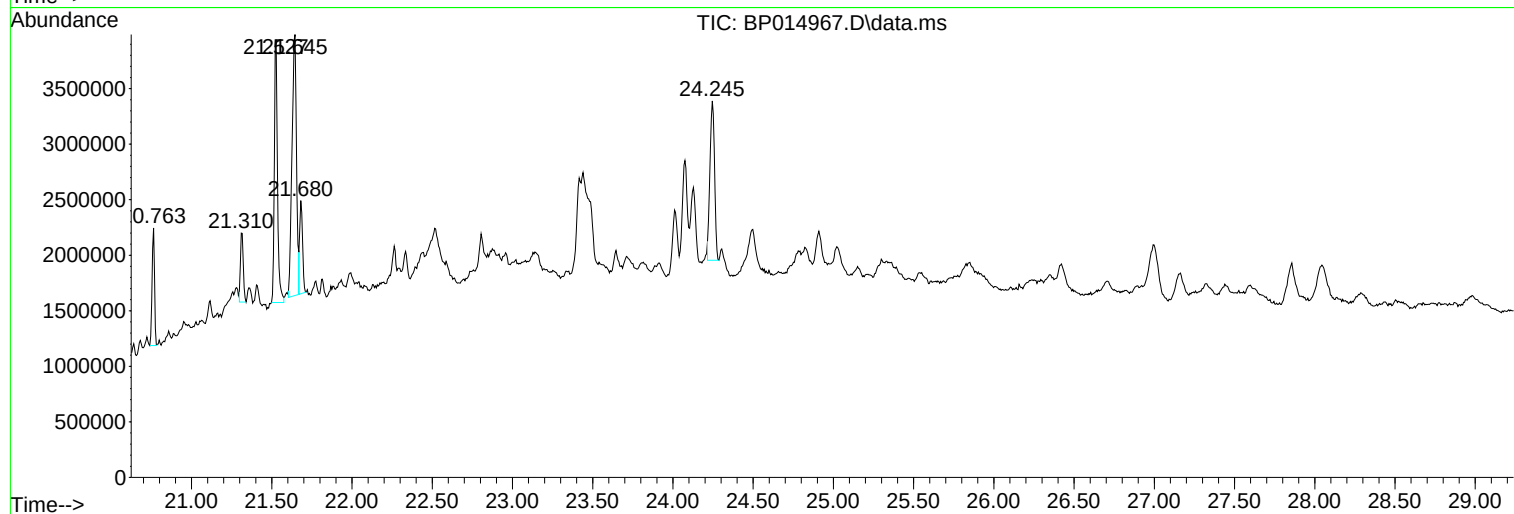
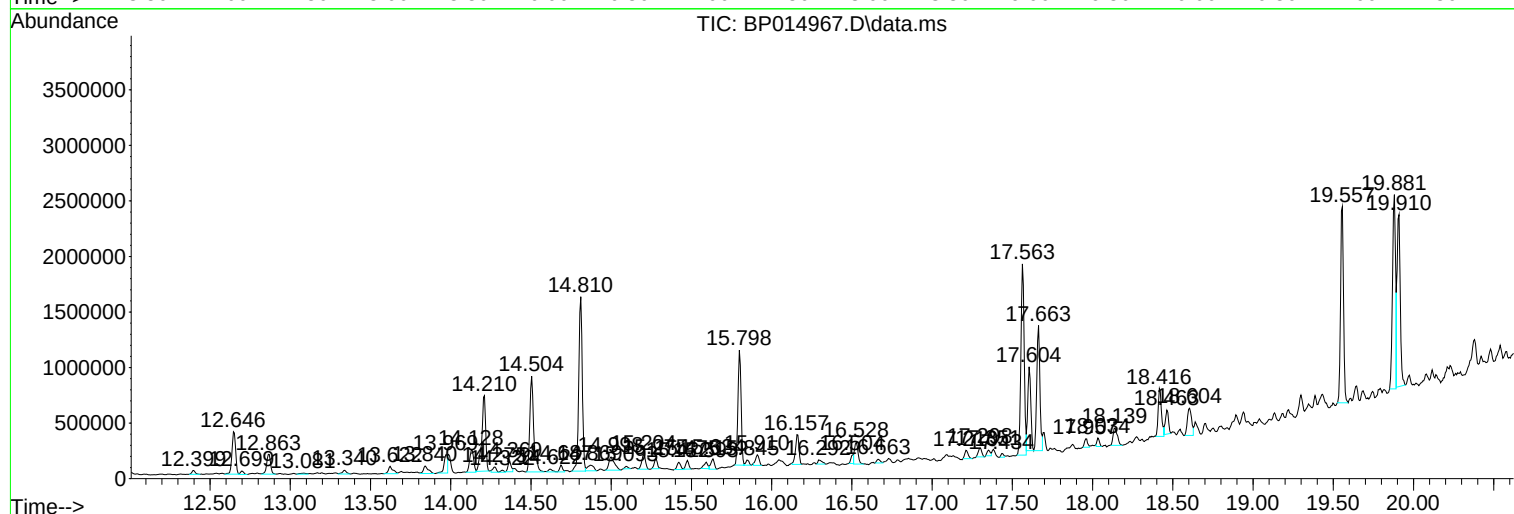
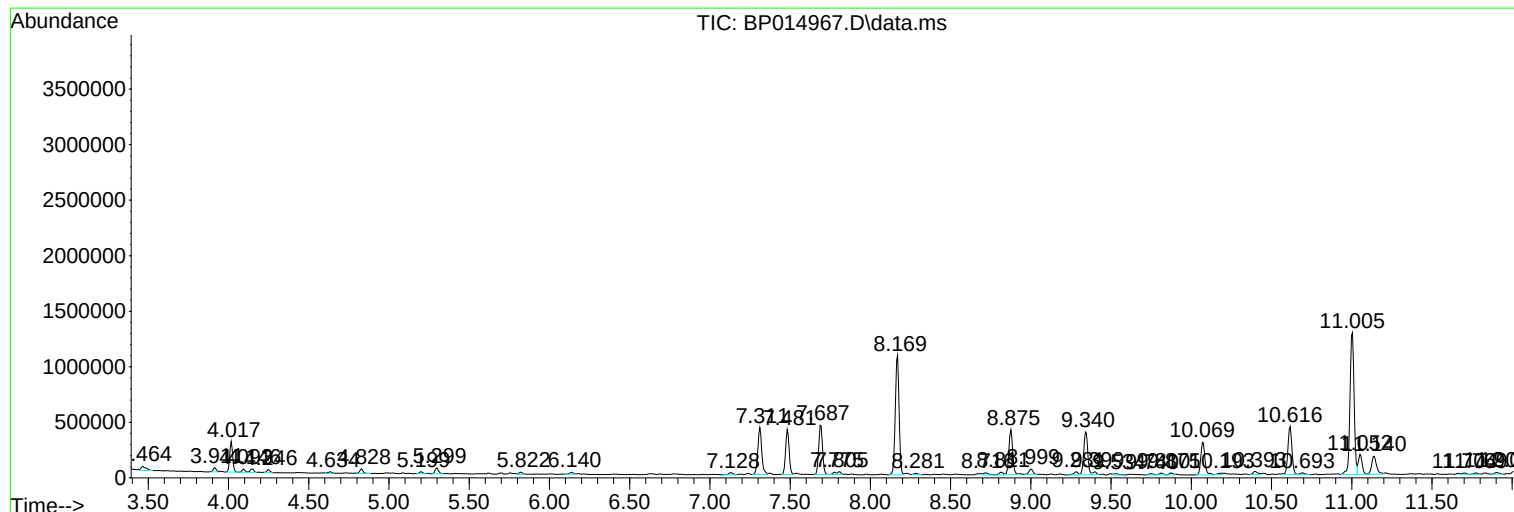
Sum of corrected areas: 49599687

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP050323.MA.M
Quant Title : SVOA CALIBRATION

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



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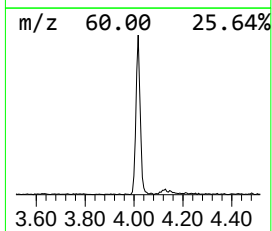
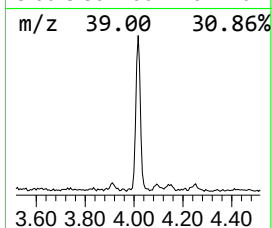
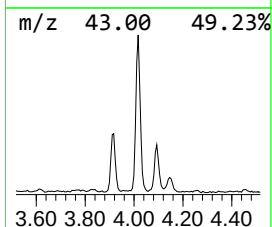
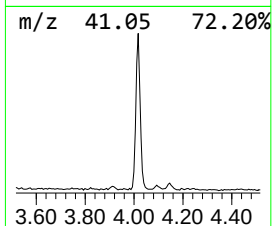
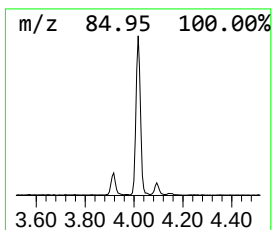
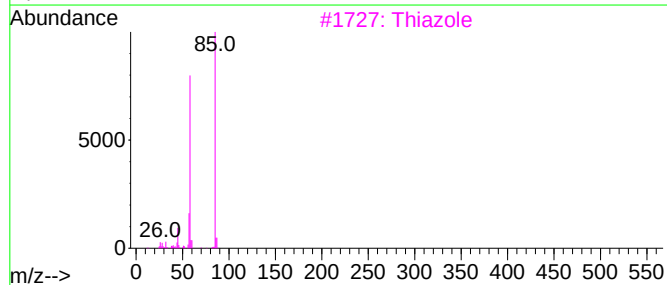
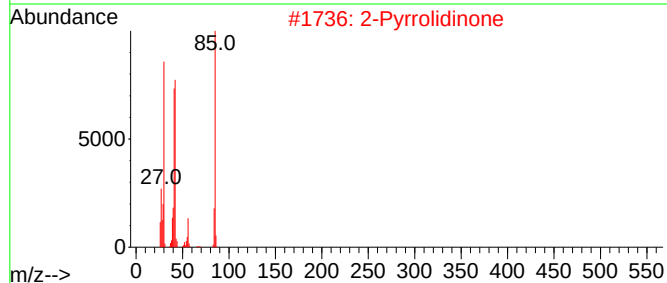
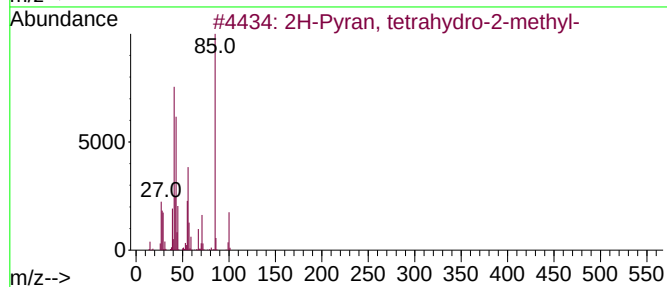
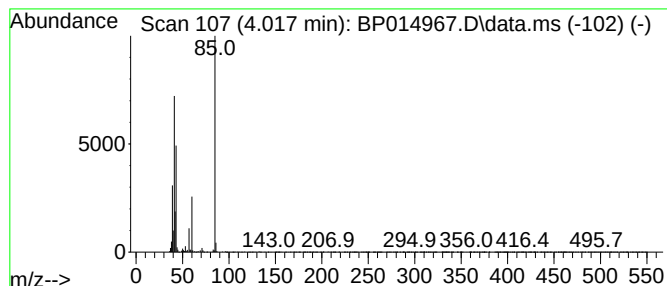
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP050323.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 unknown-01 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.017	4.03 ng/ul	350551	1,4-Dichlorobenzene-d4	8.169

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2H-Pyran, tetrahydro-2-methyl-			100	C6H12O	010141-72-7	38
2	2-Pyrrolidinone			85	C4H7NO	000616-45-5	28
3	Thiazole			85	C3H3NS	000288-47-1	28
4	2-Pyrrolidinone			85	C4H7NO	000616-45-5	9
5	2(3H)-Furanone, dihydro-5-pentyl-			156	C9H16O2	000104-61-0	9



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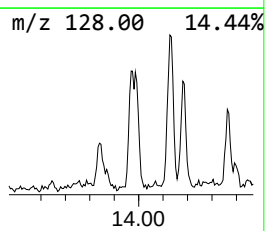
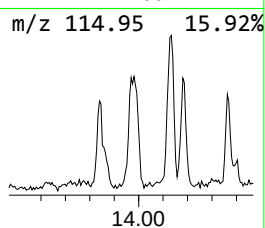
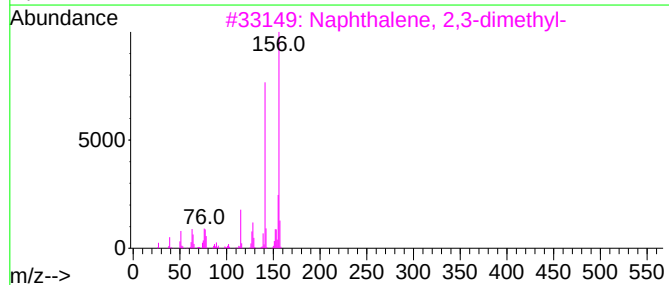
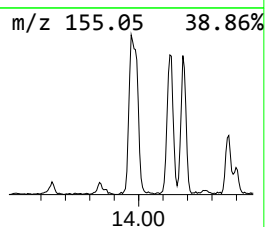
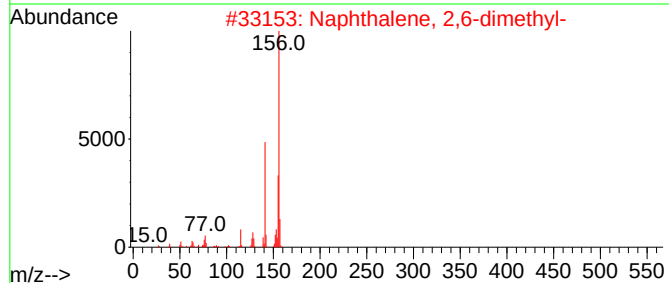
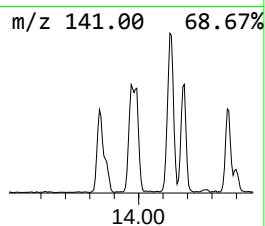
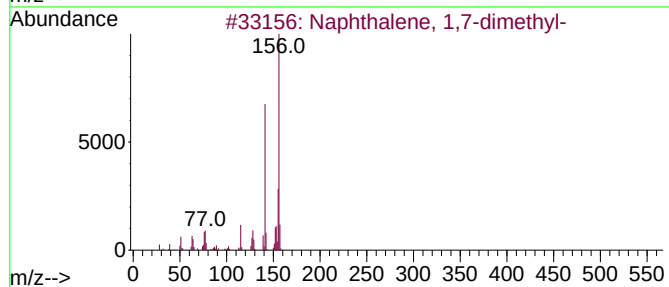
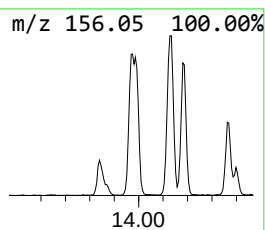
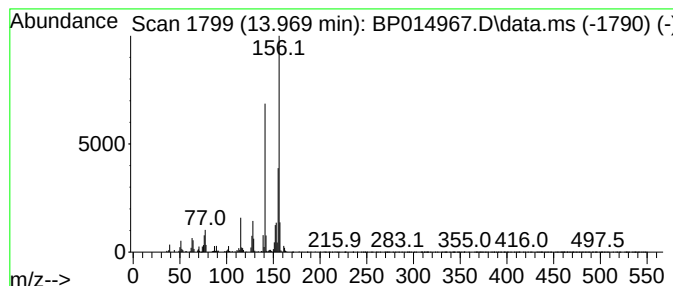
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP050323.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 Naphthalene, 1,7-dimethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.969	2.22 ng/ul	264891	Acenaphthene-d10	14.810

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1,7-dimethyl-	156	C12H12	000575-37-1	97
2			Naphthalene, 2,6-dimethyl-	156	C12H12	000581-42-0	97
3			Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	96
4			Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	96
5			Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	96



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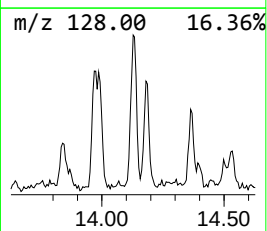
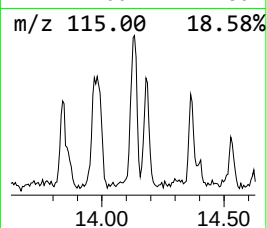
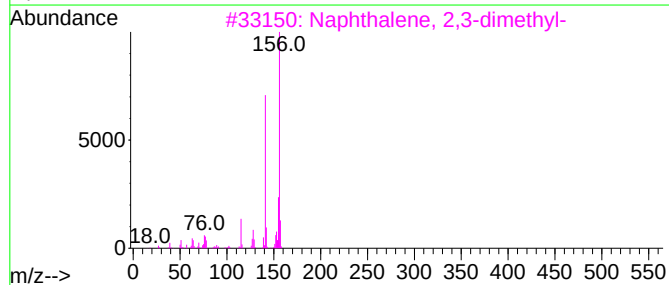
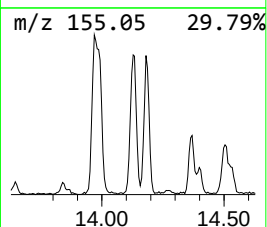
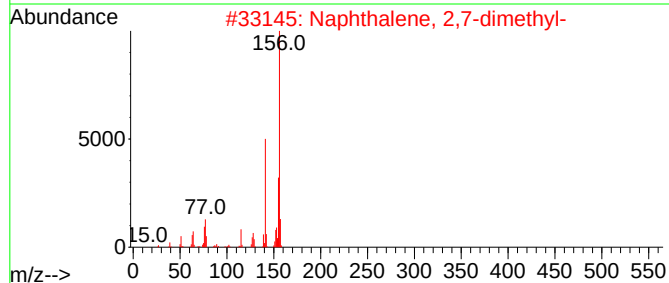
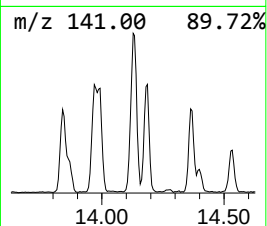
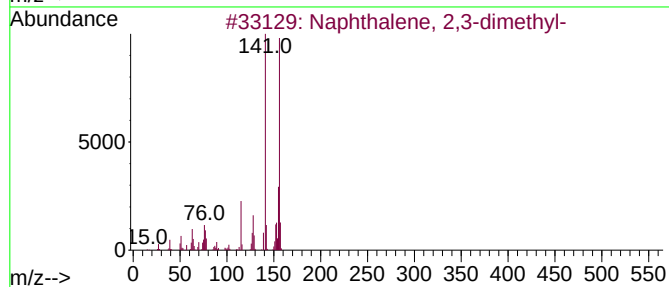
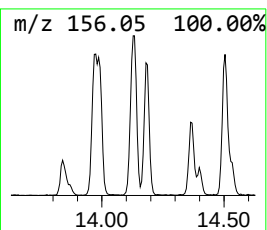
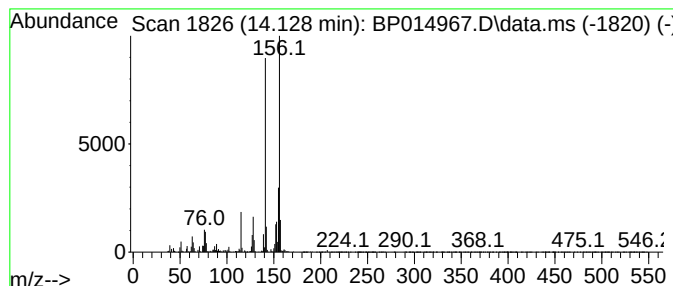
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP050323.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 Naphthalene, 2,3-dimethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.128	2.97 ng/ul	355411	Acenaphthene-d10	14.810

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	98
2		Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	97
3		Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	97
4		Naphthalene, 1,3-dimethyl-	156	C12H12	000575-41-7	97
5		Naphthalene, 1,4-dimethyl-	156	C12H12	000571-58-4	97



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP050423\
Data File : BP014967.D
Acq On : 04 May 2023 23:21
Operator : CG/JU
Sample : 02479-06
Misc :
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Instrument :
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ClientSampleId :
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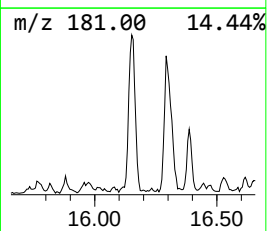
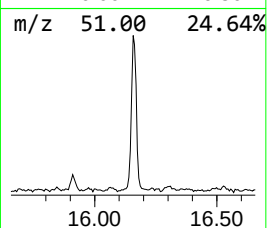
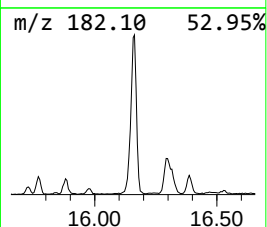
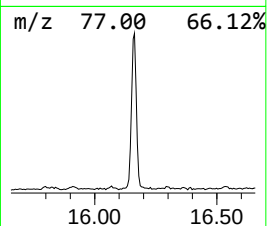
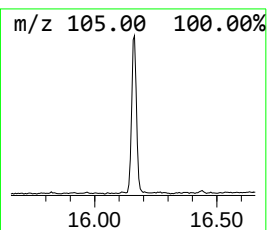
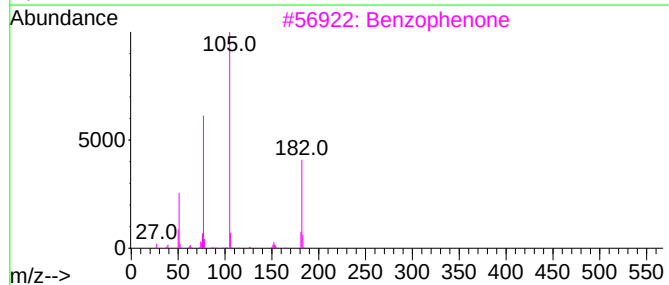
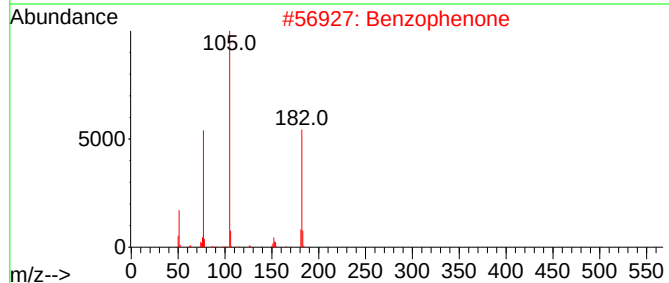
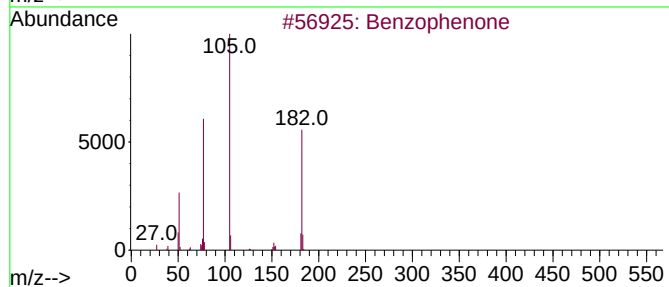
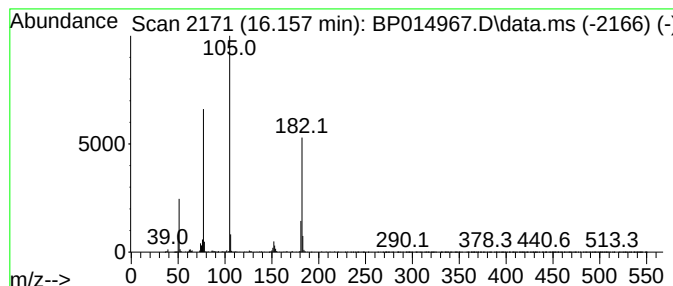
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 Benzophenone Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.157	3.35 ng/ul	400082	Acenaphthene-d10	14.810

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzophenone	182	C13H10O	000119-61-9	97
2			Benzophenone	182	C13H10O	000119-61-9	96
3			Benzophenone	182	C13H10O	000119-61-9	94
4			Benzophenone	182	C13H10O	000119-61-9	91
5			Benzophenone	182	C13H10O	000119-61-9	91



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Acq On : 04 May 2023 23:21
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Misc :
ALS Vial : 26 Sample Multiplier: 1

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ClientSampleId :
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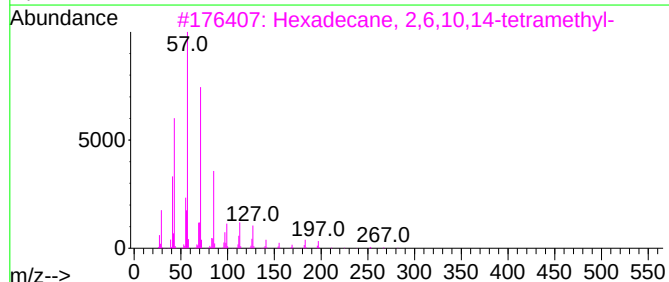
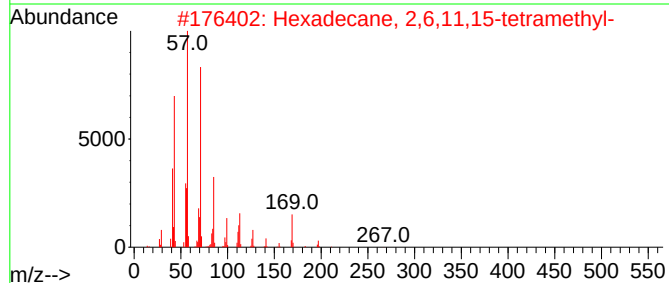
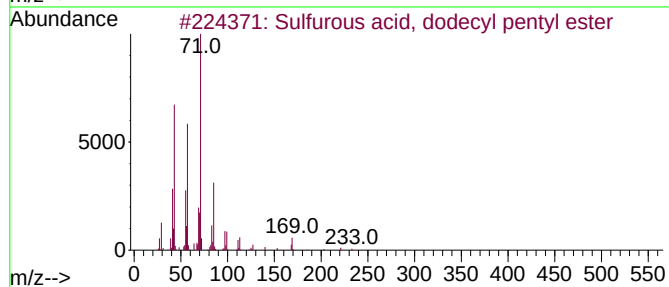
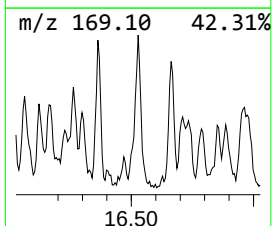
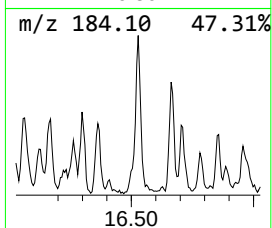
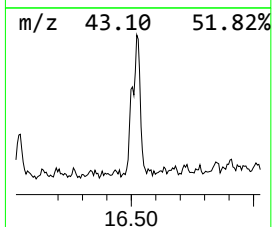
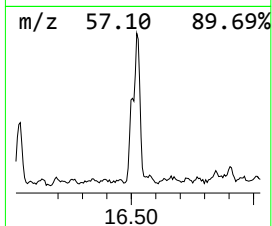
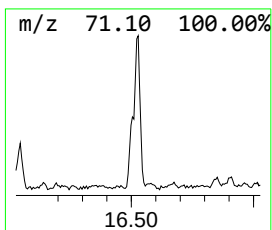
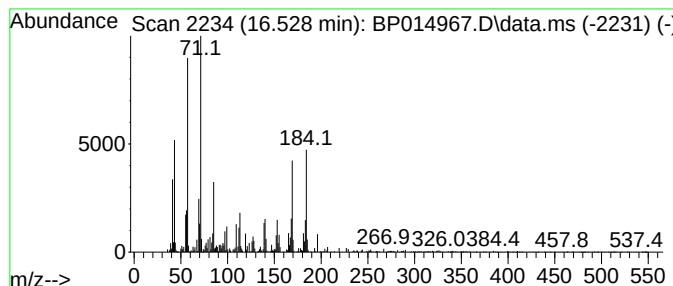
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 Sulfurous acid, dodecyl pen... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.528	2.30 ng/ul	289454	Phenanthrene-d10	17.563

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Sulfurous acid, dodecyl pentyl e...	320	C17H36O3S	1000309-14-5	50
2			Hexadecane, 2,6,11,15-tetramethyl-	282	C20H42	000504-44-9	43
3			Hexadecane, 2,6,10,14-tetramethyl-	282	C20H42	000638-36-8	38
4			Di(tetrahydrofurfuryl)adipate	314	C16H26O6	1000132-93-8	30
5			2-Hexene, 1-methoxy-, (E)-	114	C7H14O	056052-83-6	27



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP050423\
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Misc :
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Instrument :
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ClientSampleId :
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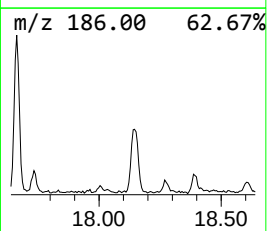
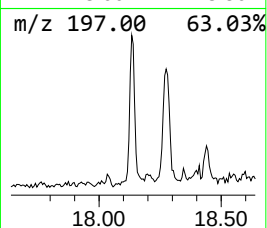
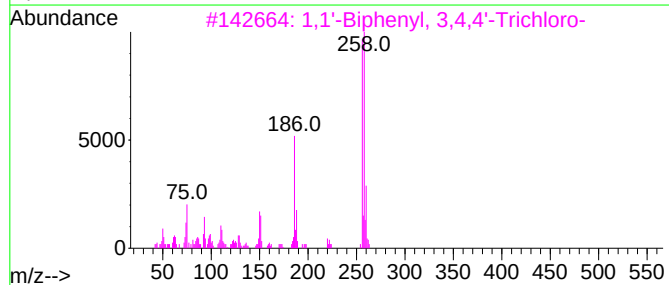
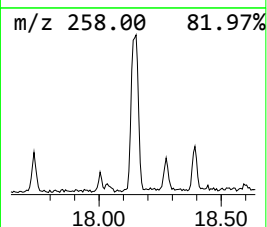
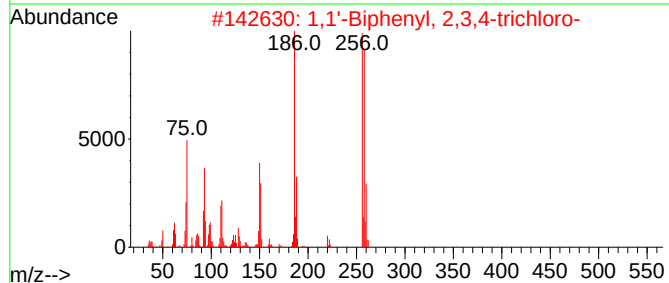
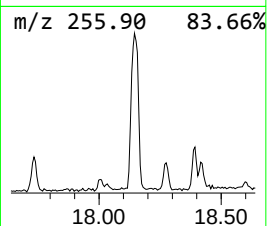
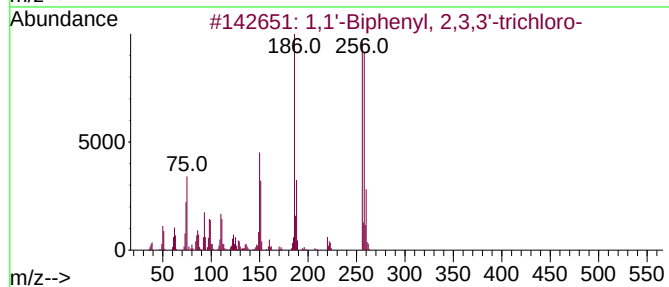
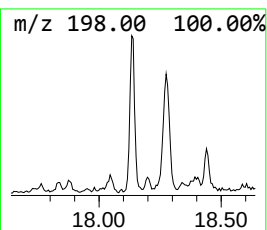
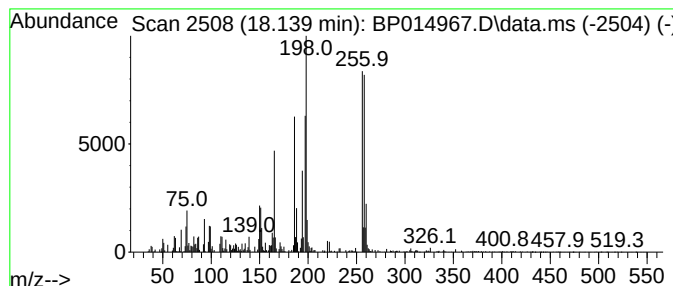
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 1,1'-Biphenyl, 2,3,3'-trich... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.139	2.30 ng/ul	290116	Phenanthrene-d10	17.563

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl, 2,3,3'-trichloro-	256	C12H7Cl3	038444-84-7	96	
2	1,1'-Biphenyl, 2,3,4-trichloro-	256	C12H7Cl3	055702-46-0	95	
3	1,1'-Biphenyl, 3,4,4'-Trichloro-	256	C12H7Cl3	038444-90-5	90	
4	1,1'-Biphenyl, 2,4',5-trichloro-	256	C12H7Cl3	016606-02-3	90	
5	2,3',6-Trichlorobiphenyl	256	C12H7Cl3	038444-76-7	87	



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP050423\
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Misc :
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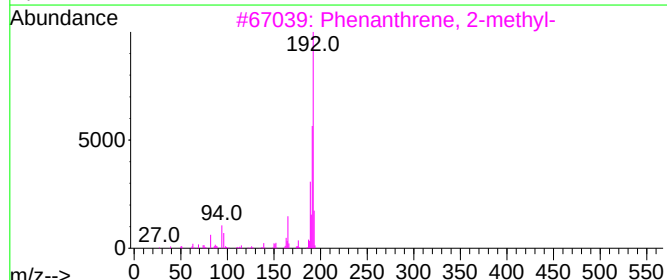
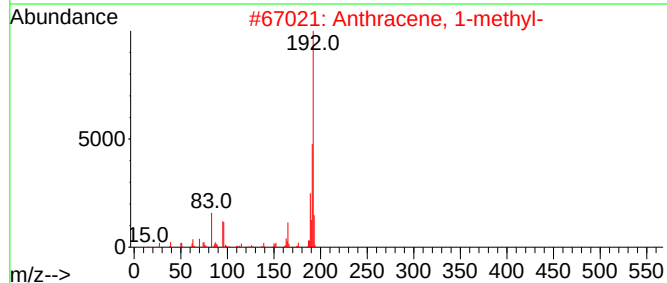
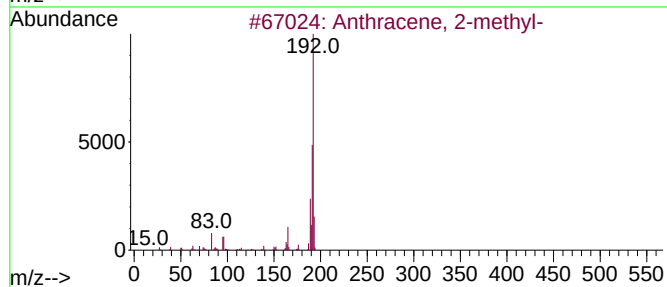
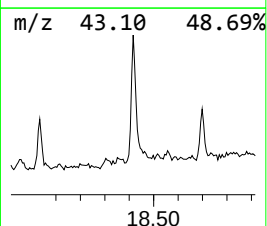
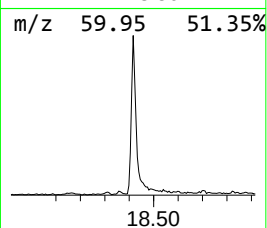
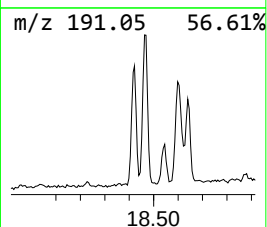
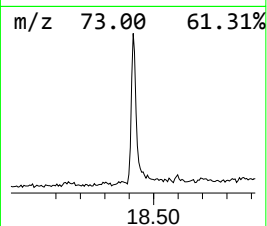
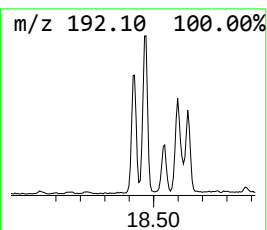
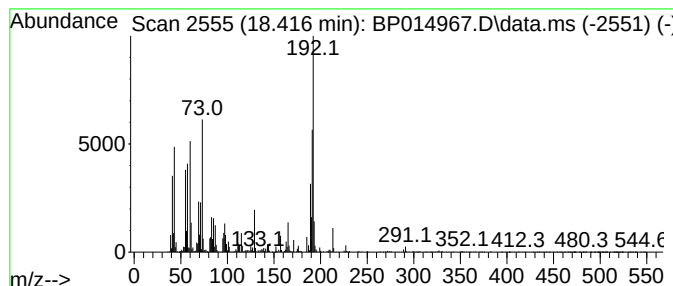
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 Anthracene, 2-methyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.416	4.98 ng/ul	626955	Phenanthrene-d10	17.563

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Anthracene, 2-methyl-	192	C15H12	000613-12-7	90
2			Anthracene, 1-methyl-	192	C15H12	000610-48-0	78
3			Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	78
4			Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	78
5			1H-Cyclopropa[1]phenanthrene,1a,...	192	C15H12	000949-41-7	78



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP050423\
Data File : BP014967.D
Acq On : 04 May 2023 23:21
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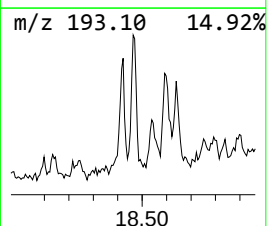
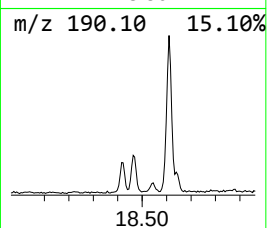
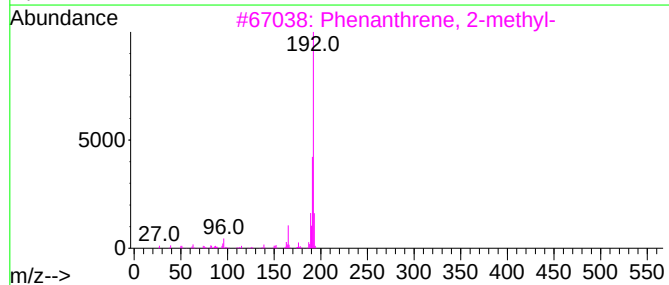
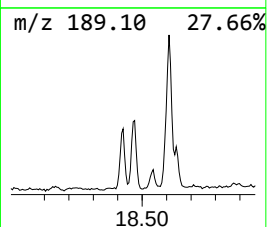
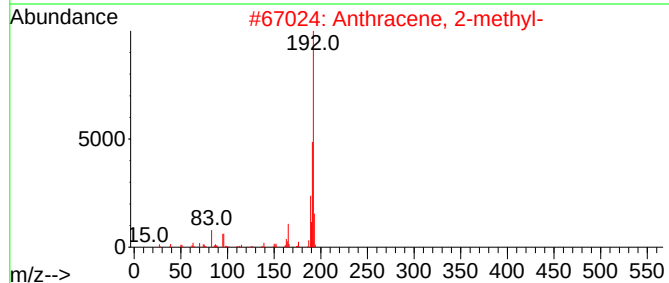
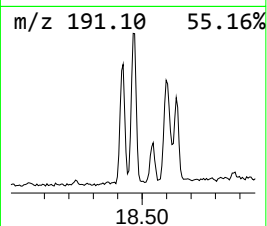
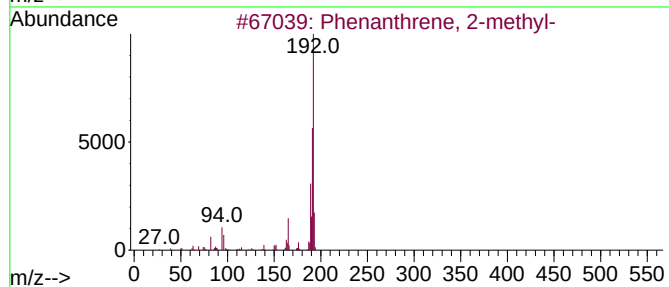
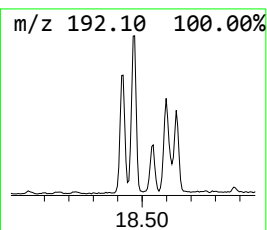
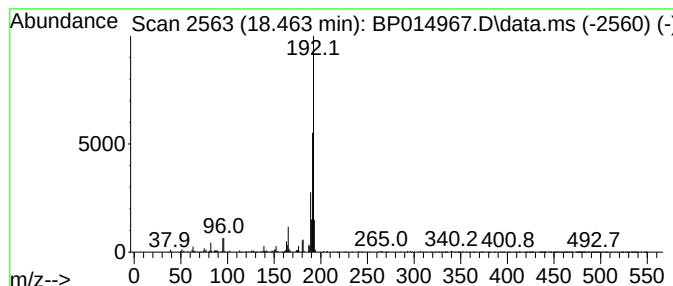
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 Phenanthrene, 2-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.463	2.24 ng/ul	282471	Phenanthrene-d10	17.563

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	97
2			Anthracene, 2-methyl-	192	C15H12	000613-12-7	96
3			Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	96
4			1H-Cyclopropa[1]phenanthrene,1a,...	192	C15H12	000949-41-7	96
5			Anthracene, 1-methyl-	192	C15H12	000610-48-0	95



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP050423\
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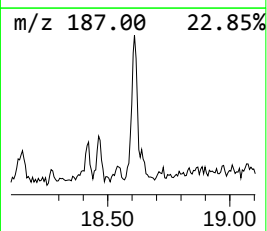
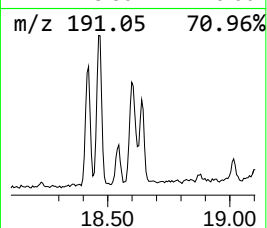
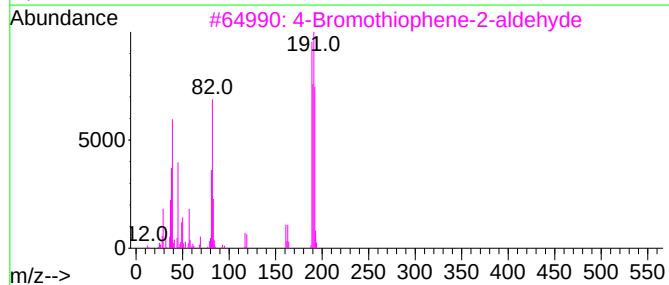
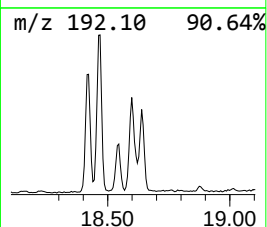
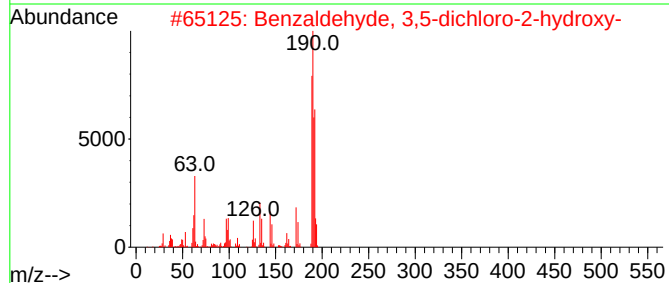
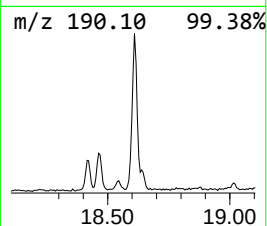
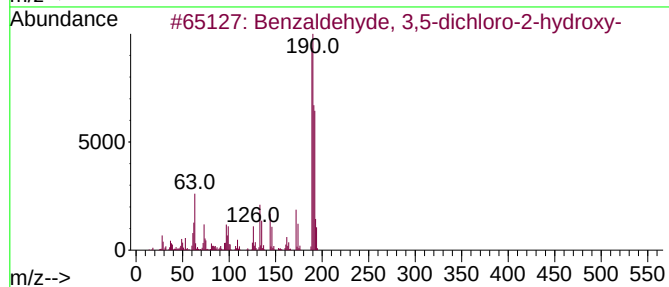
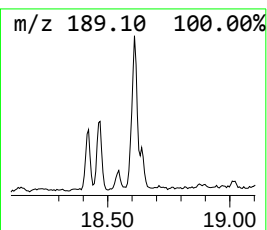
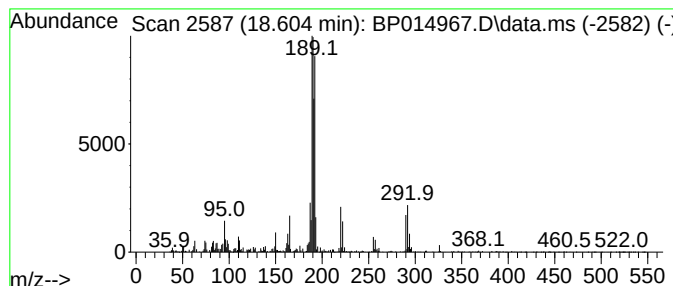
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 Benzaldehyde, 3,5-dichloro-... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.604	3.64 ng/ul	459001	Phenanthrene-d10	17.563

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzaldehyde, 3,5-dichloro-2-hyd...	190	C7H4Cl2O2	000090-60-8	64
2			Benzaldehyde, 3,5-dichloro-2-hyd...	190	C7H4Cl2O2	000090-60-8	64
3			4-Bromothiophene-2-aldehyde	190	C5H3BrOS	018791-75-8	52
4			4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	50
5			4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	49



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP050423\
Data File : BP014967.D
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ALS Vial : 26 Sample Multiplier: 1

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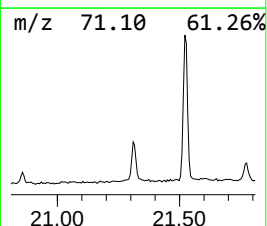
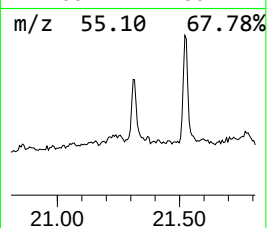
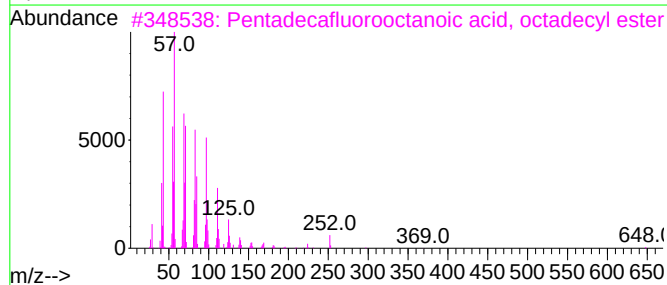
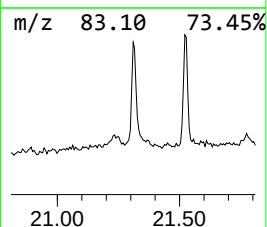
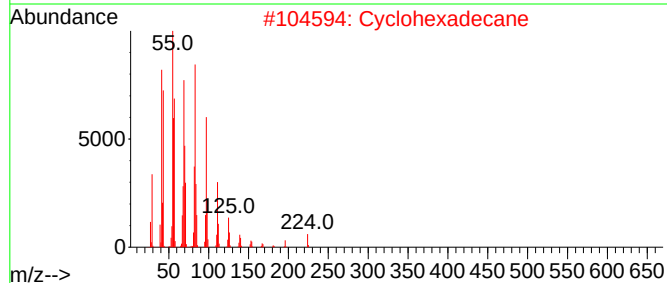
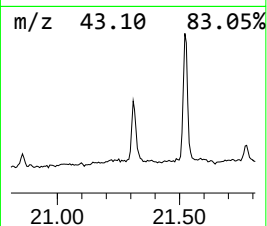
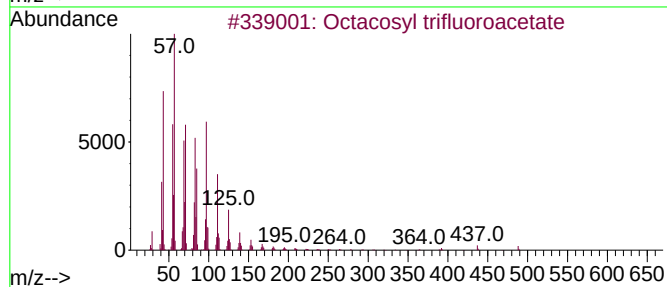
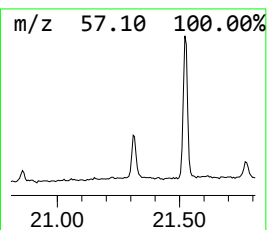
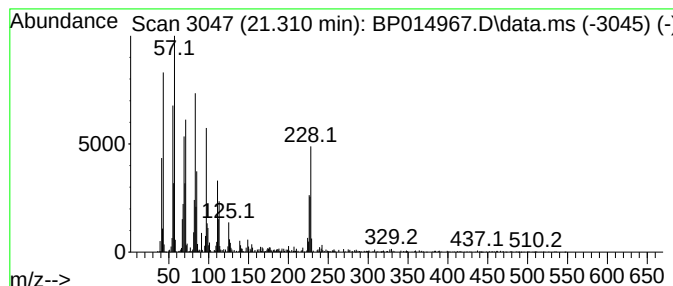
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP050323.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 10 Octacosyl trifluoroacetate Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.310	3.17 ng/ul	742030	Chrysene-d12	21.645

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octacosyl trifluoroacetate	506	C30H57F3O2	1000351-74-9	70
2			Cyclohexadecane	224	C16H32	000295-65-8	70
3			Pentadecafluorooctanoic acid, oc...	666	C26H37F15O2	1000406-04-8	70
4			Cyclopentane, (4-octyldodecyl)-	350	C25H50	005638-09-5	64
5			Pentafluoropropionic acid, octad...	416	C21H37F5O2	959261-25-7	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP050423\
Data File : BP014967.D
Acq On : 04 May 2023 23:21
Operator : CG/JU
Sample : 02479-06
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EW947

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP050323.MA.M
Quant Title : SVOA CALIBRATION

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
unknown-01	4.017	4.0	ng/ul	350551	1	8.169	1738260	20.0
Naphthalene, 1,...	13.969	2.2	ng/ul	264891	3	14.810	2391720	20.0
Naphthalene, 2,...	14.128	3.0	ng/ul	355411	3	14.810	2391720	20.0
Benzophenone	16.157	3.4	ng/ul	400082	3	14.810	2391720	20.0
Sulfurous acid,...	16.528	2.3	ng/ul	289454	4	17.563	2520050	20.0
1,1'-Biphenyl, ...	18.139	2.3	ng/ul	290116	4	17.563	2520050	20.0
Anthracene, 2-m...	18.416	5.0	ng/ul	626955	4	17.563	2520050	20.0
Phenanthrene, 2...	18.463	2.2	ng/ul	282471	4	17.563	2520050	20.0
Benzaldehyde, 3...	18.604	3.6	ng/ul	459001	4	17.563	2520050	20.0
Octacosyl trifl...	21.310	3.2	ng/ul	742030	5	21.645	4676340	20.0