

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP120123\
 Data File : BP018502.D
 Acq On : 02 Dec 2023 06:59
 Operator : MA/JU
 Sample : 05332-16DL 5X
 Misc :
 ALS Vial : 31 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 YCG97DL

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-BP120123.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BP018502.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.699	100	104	116	rBV	130920	200806	1.45%	0.183%
2	7.028	660	670	676	rVB	103898	188633	1.36%	0.172%
3	7.199	692	699	705	rBV	450584	698169	5.04%	0.636%
4	7.393	722	732	738	rBV	390875	641724	4.63%	0.585%
5	7.863	804	812	821	rBV	1622815	2546039	18.38%	2.320%
6	8.569	926	932	939	rBV	328874	513083	3.70%	0.468%
7	8.969	994	1000	1004	rVV	164960	258483	1.87%	0.236%
8	9.022	1004	1009	1020	rVB	478068	822735	5.94%	0.750%
9	9.493	1083	1089	1094	rVB	100732	153674	1.11%	0.140%
10	9.740	1122	1131	1138	rBV	301435	530057	3.83%	0.483%
11	10.057	1177	1185	1194	rBV4	133360	274634	1.98%	0.250%
12	10.157	1194	1202	1206	rVV7	34009	83203	0.60%	0.076%
13	10.281	1216	1223	1232	rVV	577600	979290	7.07%	0.892%
14	10.657	1276	1287	1293	rVV	2382386	4003257	28.90%	3.648%
15	10.710	1293	1296	1302	rVV3	60598	100694	0.73%	0.092%
16	10.798	1302	1311	1315	rVV	438596	817326	5.90%	0.745%
17	10.846	1315	1319	1326	rVV2	405406	790369	5.71%	0.720%
18	10.904	1326	1329	1333	rVV	95999	161136	1.16%	0.147%
19	10.957	1333	1338	1344	rVB	139683	242064	1.75%	0.221%
20	11.075	1353	1358	1365	rVV	233514	461599	3.33%	0.421%
21	11.140	1365	1369	1374	rVB2	60290	90868	0.66%	0.083%
22	11.246	1381	1387	1393	rBV2	55435	98974	0.71%	0.090%
23	11.322	1393	1400	1404	rVB3	54014	93387	0.67%	0.085%
24	11.428	1412	1418	1423	rBV	78253	134660	0.97%	0.123%
25	11.575	1438	1443	1446	rBV	56594	92141	0.67%	0.084%
26	11.704	1460	1465	1469	rVV	88760	158671	1.15%	0.145%
27	11.763	1469	1475	1479	rVV6	70857	168645	1.22%	0.154%
28	11.851	1485	1490	1498	rVB	114945	201832	1.46%	0.184%
29	12.045	1519	1523	1529	rVB2	81857	119189	0.86%	0.109%
30	12.210	1547	1551	1556	rBV	207668	325288	2.35%	0.296%
31	12.375	1573	1579	1584	rBV	119356	191110	1.38%	0.174%
32	12.434	1584	1589	1594	rVV5	49510	103555	0.75%	0.094%
33	12.504	1594	1601	1614	rVB4	137162	406488	2.93%	0.370%
34	13.328	1734	1741	1747	rVV	2908037	4133682	29.84%	3.767%
35	13.398	1747	1753	1761	rVV	166019	291505	2.10%	0.266%
36	13.504	1767	1771	1775	rVV4	58263	99882	0.72%	0.091%

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Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120123.MA.M
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37	13.551	1775	1779	1786	rVB	134556	196444	1.42%	0.179%
38	13.669	1794	1799	1806	rVV2	65093	145468	1.05%	0.133%
39	13.816	1815	1824	1830	rBV	849464	1264982	9.13%	1.153%
40	13.904	1830	1839	1851	rVB	1318793	1957435	14.13%	1.784%
41	14.051	1851	1864	1867	rBV	532411	906401	6.54%	0.826%
42	14.081	1867	1869	1875	rVB	297542	348899	2.52%	0.318%
43	14.216	1880	1892	1898	rBV2	8031337	13852696	100.00%	12.625%
44	14.492	1928	1939	1944	rBV	3507650	5342849	38.57%	4.869%
45	14.551	1944	1949	1959	rVB	4408643	6718714	48.50%	6.123%
46	14.687	1966	1972	1982	rBV3	87570	210415	1.52%	0.192%
47	14.887	2000	2006	2014	rVB	3085825	4527979	32.69%	4.127%
48	14.963	2015	2019	2029	rVB3	77923	142237	1.03%	0.130%
49	15.087	2034	2040	2048	rBV3	168008	407526	2.94%	0.371%
50	15.169	2048	2054	2064	rVB3	135378	267684	1.93%	0.244%
51	15.257	2064	2069	2075	rBV2	133213	283631	2.05%	0.258%
52	15.310	2075	2078	2083	rVB2	91365	131575	0.95%	0.120%
53	15.439	2093	2100	2102	rBV	194260	281146	2.03%	0.256%
54	15.481	2102	2107	2111	rVV	1922360	2682740	19.37%	2.445%
55	15.534	2111	2116	2122	rVV	3713397	5541402	40.00%	5.050%
56	15.592	2122	2126	2130	rVV2	322514	457427	3.30%	0.417%
57	15.628	2130	2132	2135	rVV	94846	126162	0.91%	0.115%
58	15.663	2135	2138	2140	rVV	96543	119297	0.86%	0.109%
59	15.698	2140	2144	2148	rVV3	81100	125137	0.90%	0.114%
60	15.745	2148	2152	2155	rVV2	67748	97108	0.70%	0.088%
61	15.786	2155	2159	2162	rVV	173838	223970	1.62%	0.204%
62	15.828	2162	2166	2171	rVB	295878	406552	2.93%	0.371%
63	15.975	2186	2191	2199	rVB	390797	707759	5.11%	0.645%
64	16.069	2199	2207	2213	rVB3	67055	148002	1.07%	0.135%
65	16.210	2225	2231	2237	rBV	222633	324303	2.34%	0.296%
66	16.498	2275	2280	2282	rBV2	87556	135658	0.98%	0.124%
67	16.539	2283	2287	2292	rVV	137873	232703	1.68%	0.212%
68	16.886	2341	2346	2353	rVB	412547	605584	4.37%	0.552%
69	17.051	2370	2374	2380	rVV	227125	327425	2.36%	0.298%
70	17.239	2396	2406	2410	rBV	3965385	5847942	42.22%	5.330%
71	17.281	2410	2413	2417	rVV	376794	497820	3.59%	0.454%
72	17.333	2417	2422	2426	rVV2	2244410	3397073	24.52%	3.096%
73	17.369	2426	2428	2433	rVB	669949	780339	5.63%	0.711%
74	17.439	2433	2440	2448	rVB	1285825	1815580	13.11%	1.655%
75	17.639	2465	2474	2479	rBV	972736	1449985	10.47%	1.321%
76	17.798	2498	2501	2510	rVB2	138964	215233	1.55%	0.196%

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 Title : SVOA CALIBRATION

77	18.069	2543	2547	2549	rBV	84815	101106	0.73%	0.092%
78	18.104	2549	2553	2555	rVV	132914	178221	1.29%	0.162%
79	18.145	2555	2560	2563	rVV2	289049	537754	3.88%	0.490%
80	18.186	2563	2567	2570	rVV	442602	565155	4.08%	0.515%
81	18.222	2570	2573	2578	rVB2	168768	225834	1.63%	0.206%
82	18.292	2579	2585	2589	rBV	459241	703765	5.08%	0.641%
83	18.328	2589	2591	2596	rVB	107835	121724	0.88%	0.111%
84	18.398	2596	2603	2606	rBV4	165195	282755	2.04%	0.258%
85	18.439	2606	2610	2613	rVB	210406	257278	1.86%	0.234%
86	18.528	2620	2625	2631	rBV	321463	457618	3.30%	0.417%
87	18.586	2631	2635	2638	rVV2	115227	166699	1.20%	0.152%
88	18.622	2638	2641	2646	rVB	184546	235644	1.70%	0.215%
89	19.127	2723	2727	2734	rVB3	86816	157371	1.14%	0.143%
90	19.245	2742	2747	2752	rVB	1051364	1321537	9.54%	1.204%
91	19.327	2758	2761	2764	rBV2	77119	100652	0.73%	0.092%
92	19.363	2764	2767	2770	rVB2	96814	95052	0.69%	0.087%
93	19.569	2796	2802	2805	rBV	2343620	3230483	23.32%	2.944%
94	19.598	2805	2807	2816	rVB	959080	1052897	7.60%	0.960%
95	20.027	2875	2880	2886	rBV	503650	713925	5.15%	0.651%
96	20.651	2982	2986	2992	rBV2	164645	222872	1.61%	0.203%
97	21.321	3094	3100	3105	rBV	4413510	5523413	39.87%	5.034%
98	23.539	3471	3477	3483	rBV	1458557	2640148	19.06%	2.406%
99	23.604	3485	3488	3495	rVB	280688	454403	3.28%	0.414%
100	23.692	3497	3503	3516	rVB	3042317	5952201	42.97%	5.425%

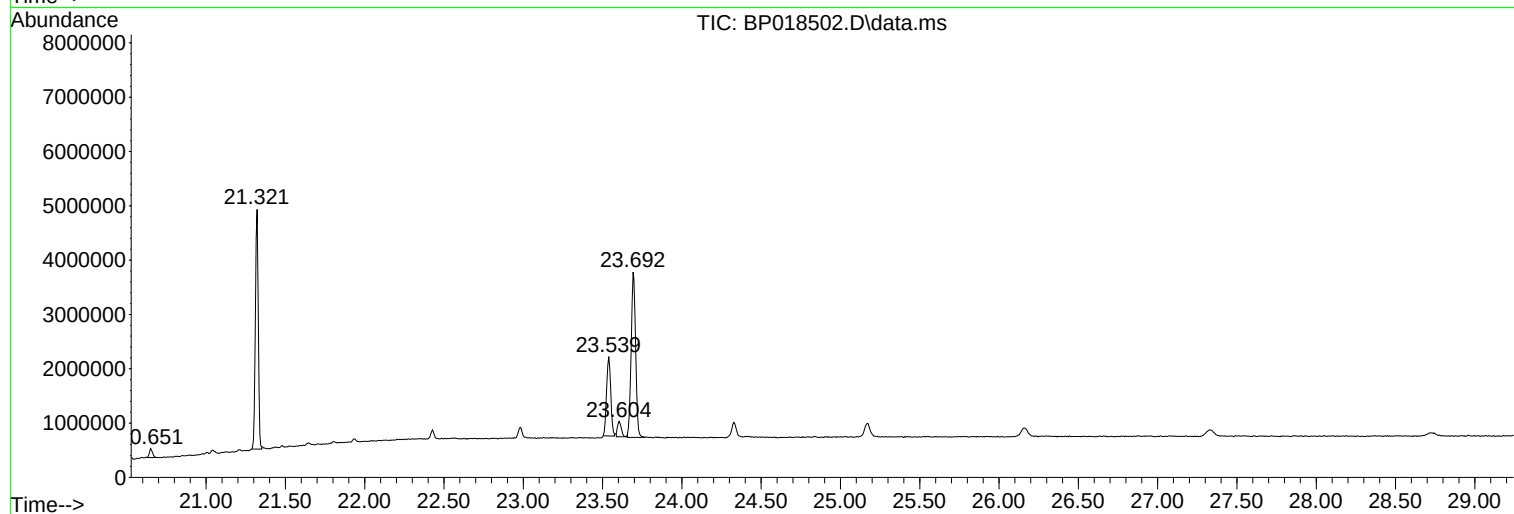
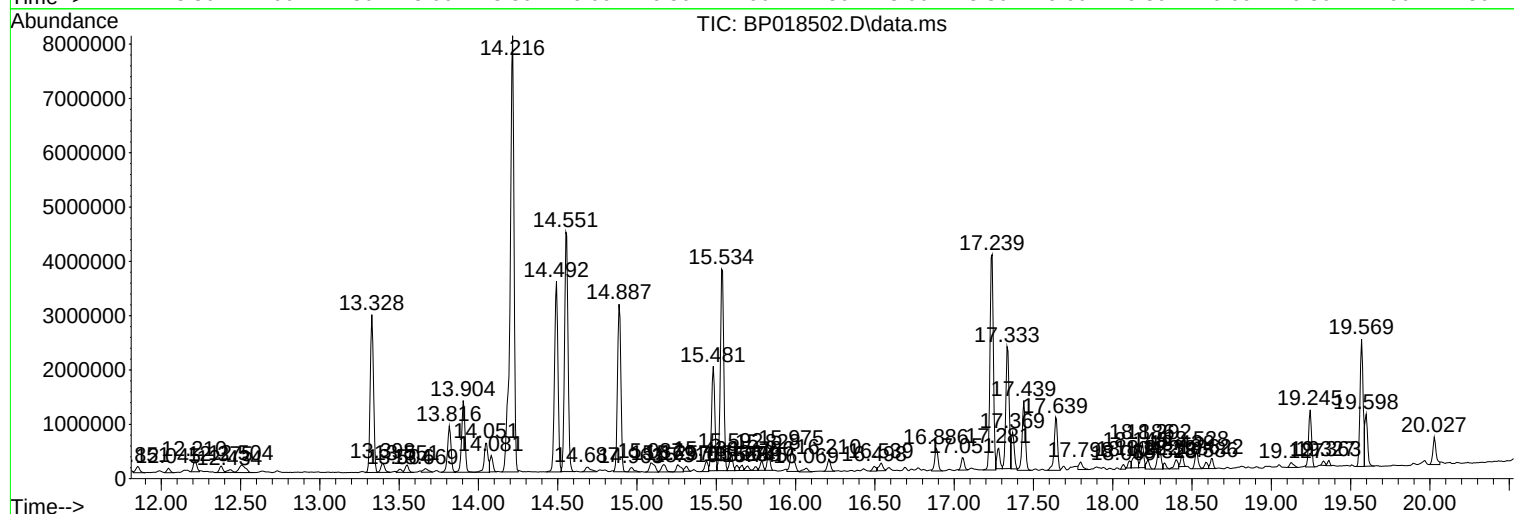
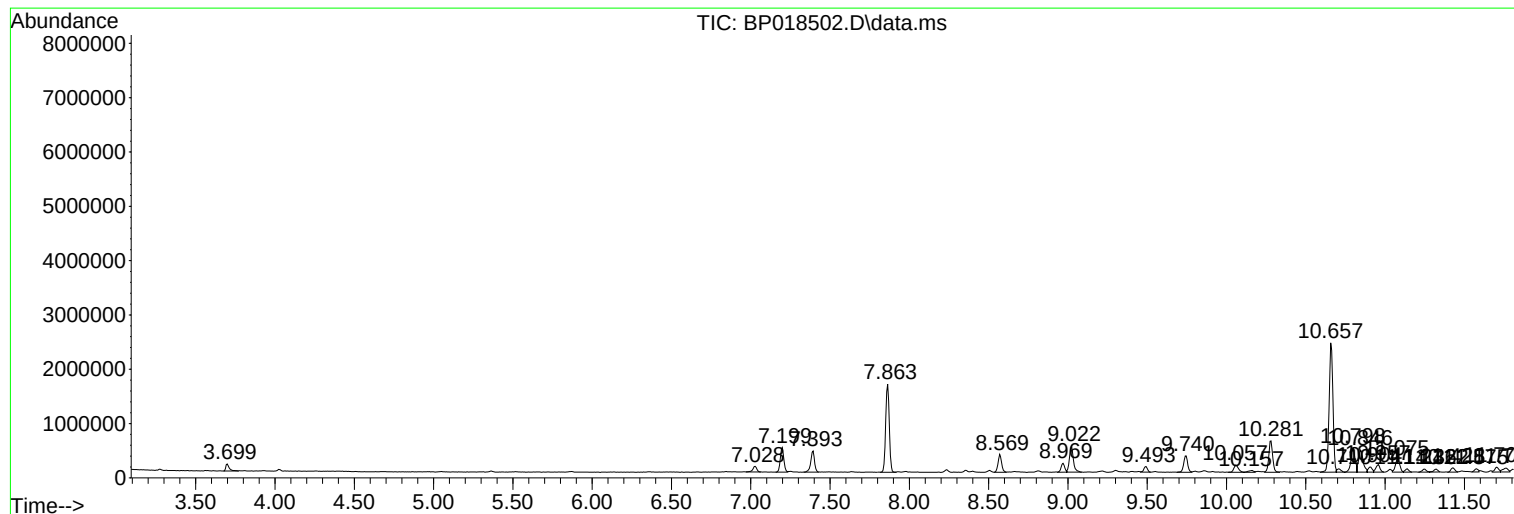
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120123.MA.M
Quant Title : SVOA CALIBRATION

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



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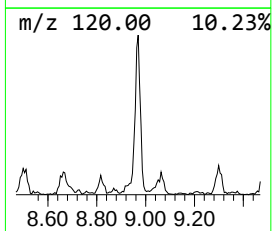
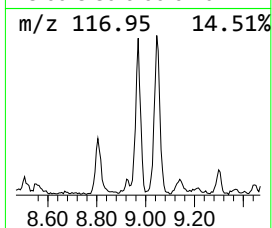
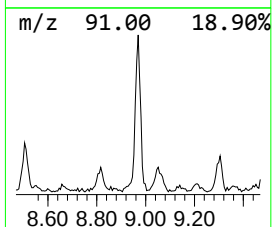
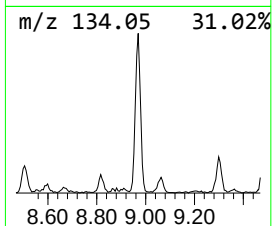
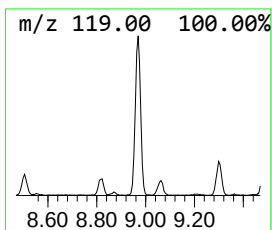
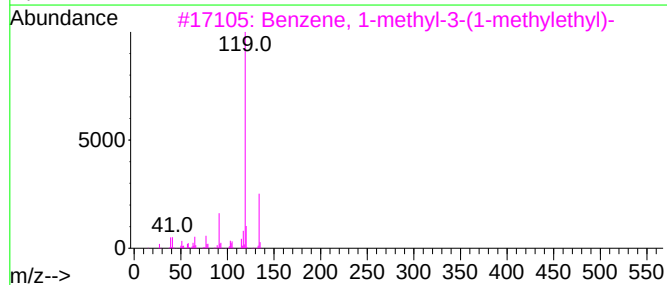
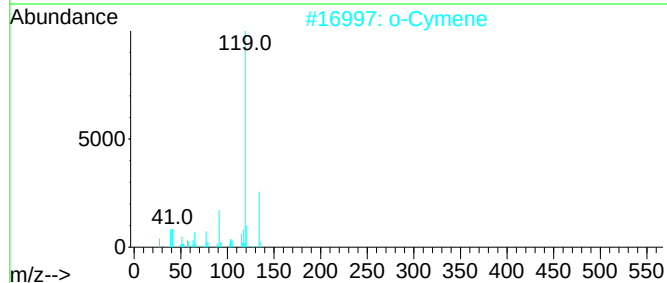
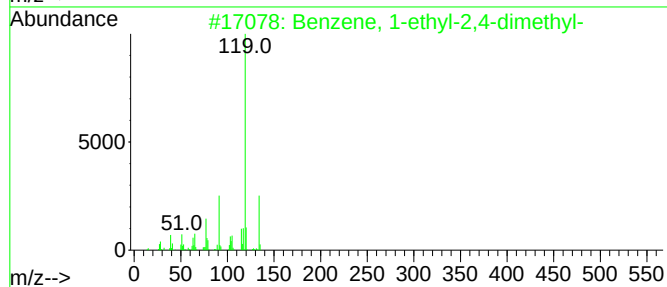
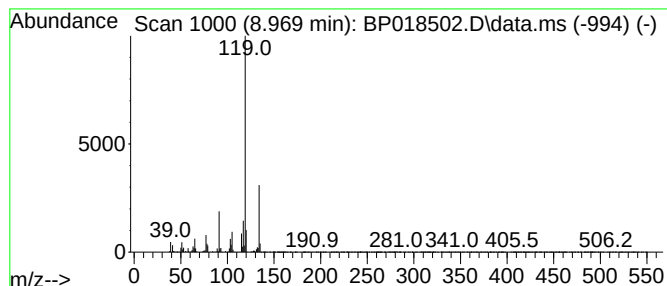
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120123.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Benzene, 1-ethyl-2,4-dimethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.969	2.03 ng/ul	258483	1,4-Dichlorobenzene-d4	7.863

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	95
2			o-Cymene	134	C10H14	000527-84-4	95
3			Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	94
4			Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	94
5			Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	93



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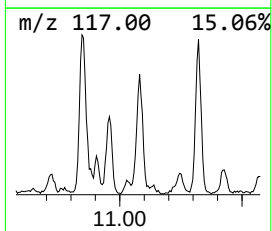
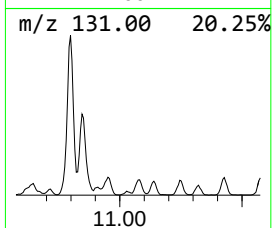
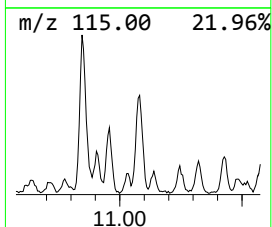
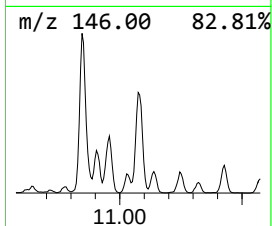
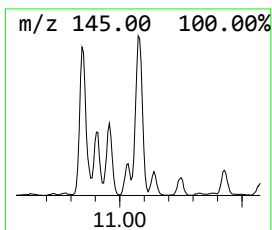
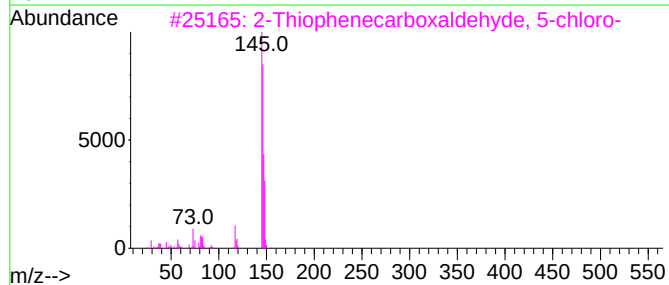
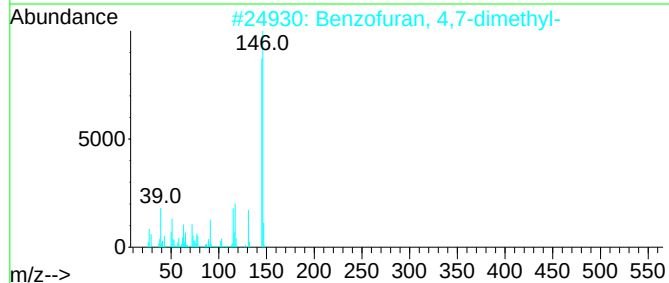
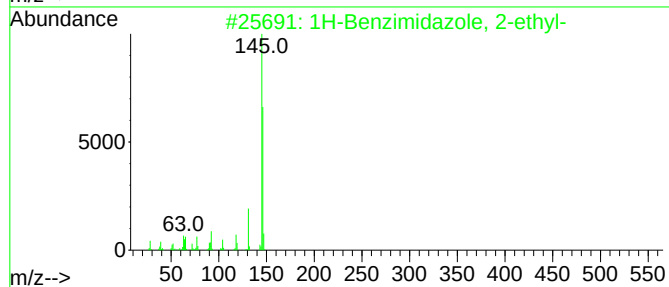
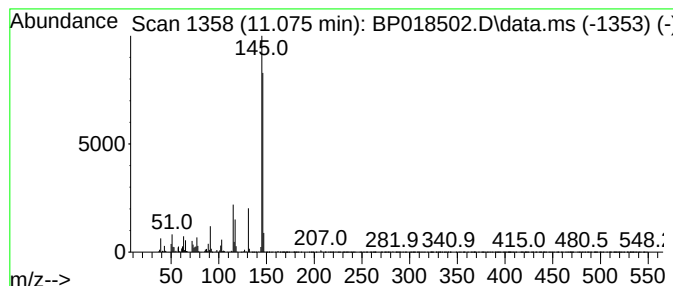
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120123.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 1H-Benzimidazole, 2-ethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.075	2.31 ng/ul	461599	Naphthalene-d8	10.657

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1H-Benzimidazole, 2-ethyl-	146	C9H10N2	001848-84-6	64
2			Benzofuran, 4,7-dimethyl-	146	C10H10O	028715-26-6	58
3			2-Thiophenecarboxaldehyde, 5-chl...	146	C5H3ClOS	007283-96-7	53
4			1H-1,3-Benzodiazole-5-carbaldehyde	146	C8H6N2O	058442-17-4	53
5			1H-Benzimidazole, 5,6-dimethyl-	146	C9H10N2	000582-60-5	50



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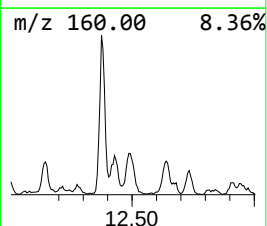
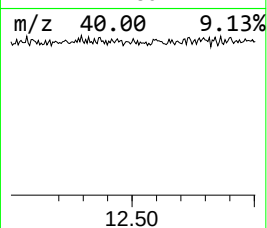
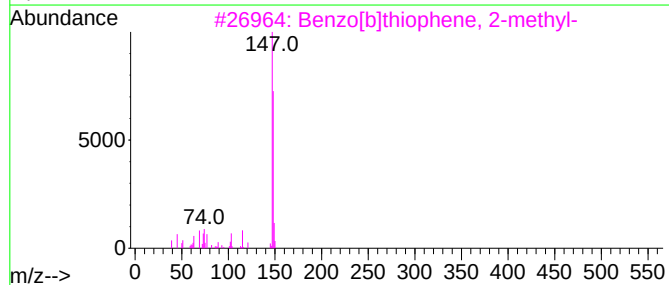
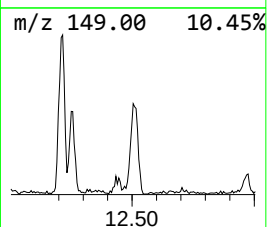
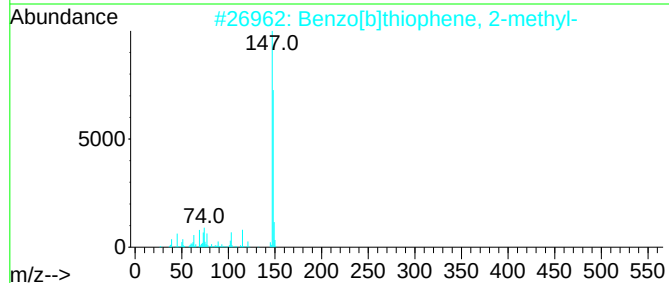
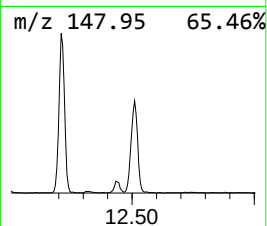
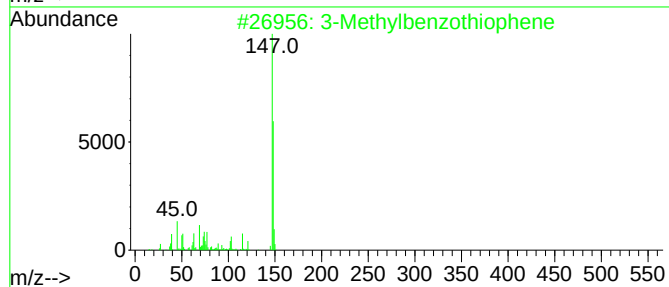
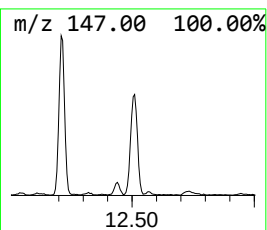
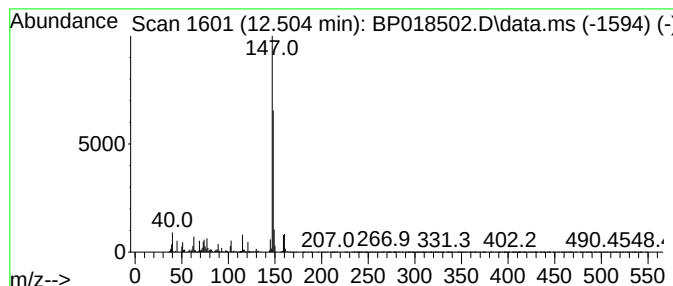
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Quant Title : SVOA CALIBRATION

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

Peak Number 3 3-Methylbenzothiophene Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.504	2.03 ng/ul	406488	Naphthalene-d8	10.657

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3-Methylbenzothiophene	148	C9H8S	001455-18-1	90
2			Benzo[b]thiophene, 2-methyl-	148	C9H8S	001195-14-8	90
3			Benzo[b]thiophene, 2-methyl-	148	C9H8S	001195-14-8	90
4			3-Methylbenzothiophene	148	C9H8S	001455-18-1	87
5			Benzo[b]thiophene, 4-methyl-	148	C9H8S	014315-11-8	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP120123\
Data File : BP018502.D
Acq On : 02 Dec 2023 06:59
Operator : MA/JU
Sample : 05332-16DL 5X
Misc :
ALS Vial : 31 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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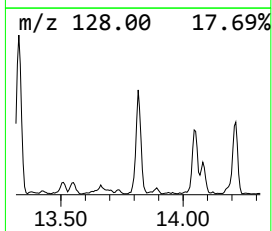
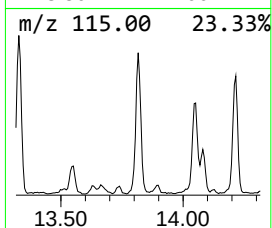
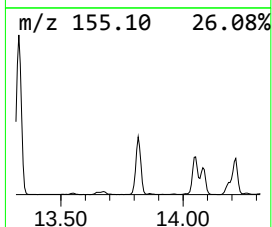
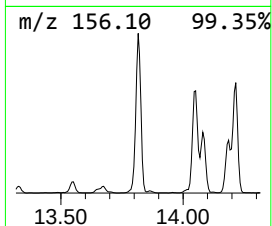
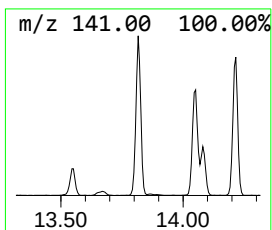
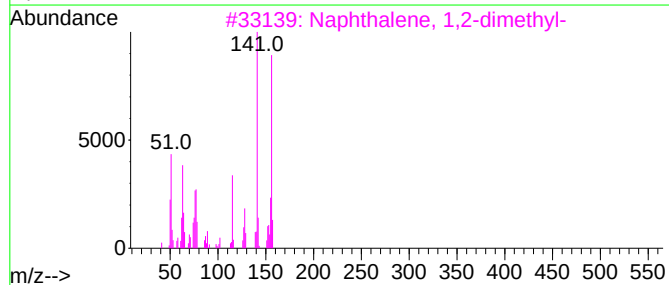
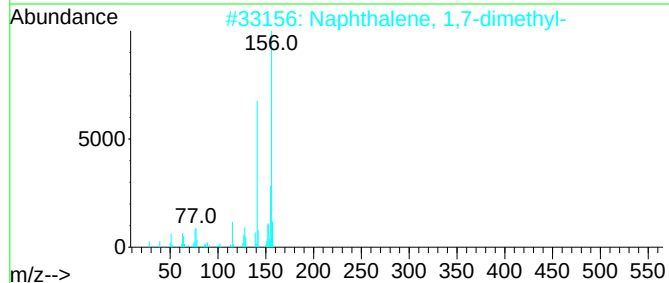
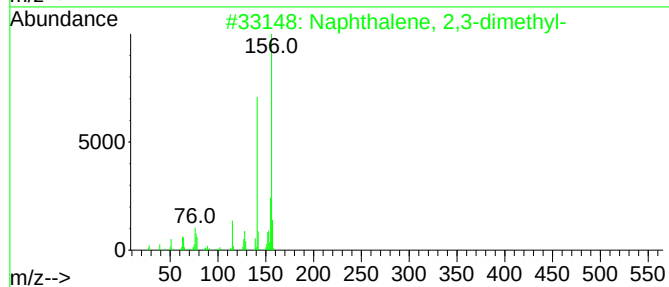
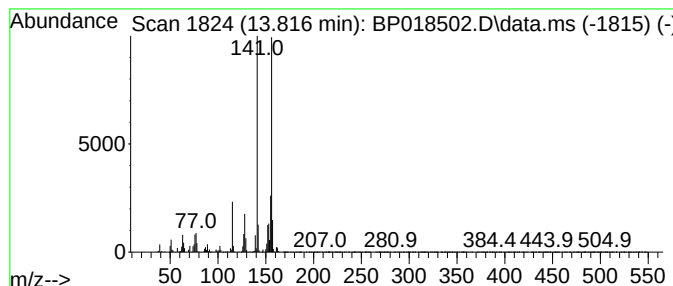
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 Naphthalene, 2,3-dimethyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.816	4.74 ng/ul	1264980	Acenaphthene-d10	14.492

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	97
2			Naphthalene, 1,7-dimethyl-	156	C12H12	000575-37-1	97
3			Naphthalene, 1,2-dimethyl-	156	C12H12	000573-98-8	97
4			Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	96
5			Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	96



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP120123\
Data File : BP018502.D
Acq On : 02 Dec 2023 06:59
Operator : MA/JU
Sample : 05332-16DL 5X
Misc :
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Instrument :
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ClientSampleId :
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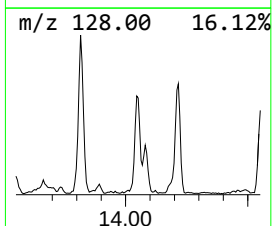
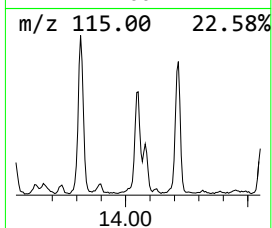
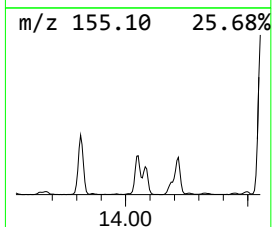
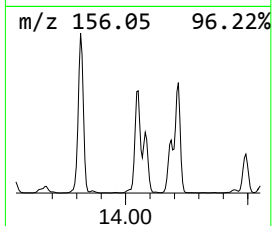
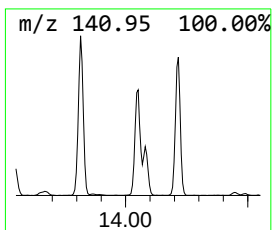
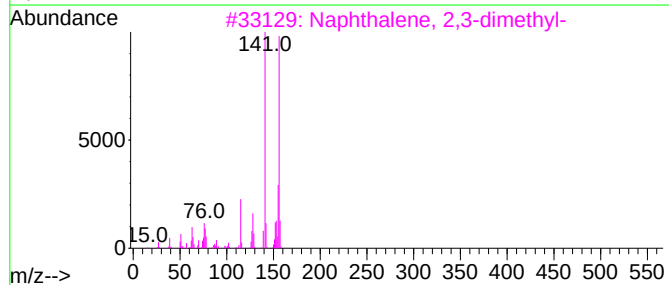
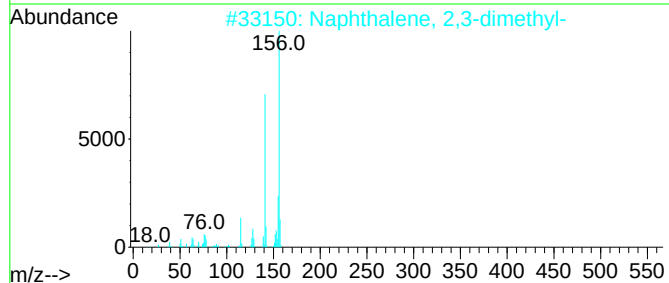
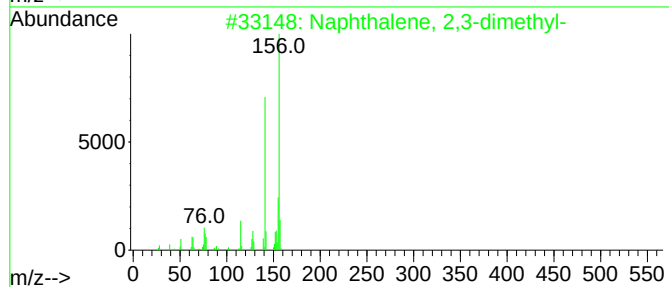
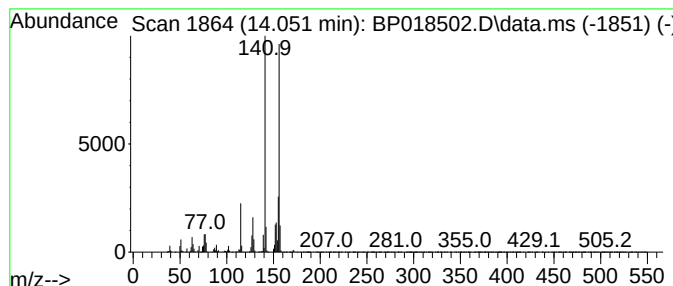
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 Naphthalene, 1,7-dimethyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.051	3.39 ng/ul	906401	Acenaphthene-d10	14.492

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



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP120123\
Data File : BP018502.D
Acq On : 02 Dec 2023 06:59
Operator : MA/JU
Sample : 05332-16DL 5X
Misc :
ALS Vial : 31 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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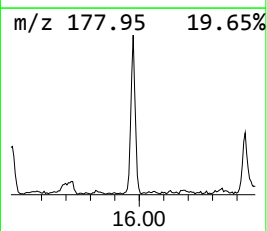
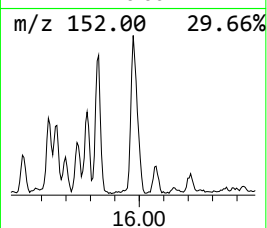
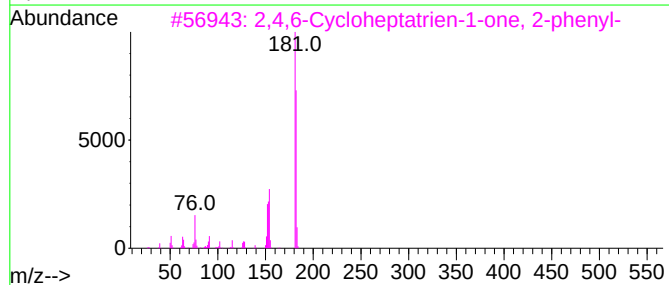
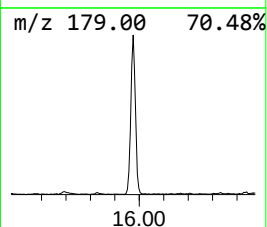
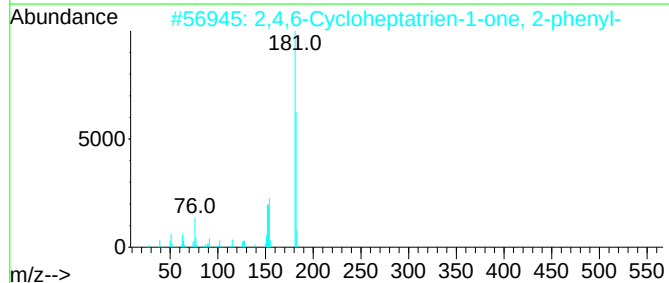
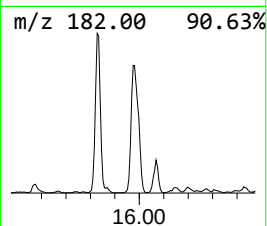
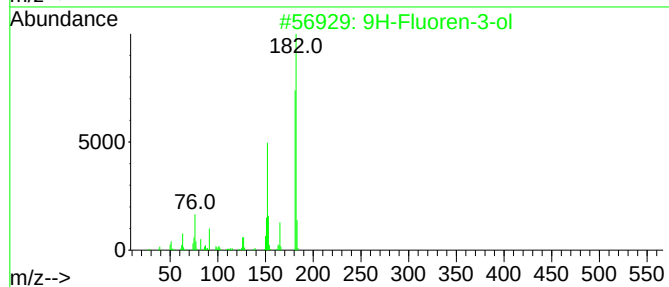
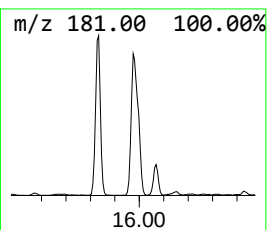
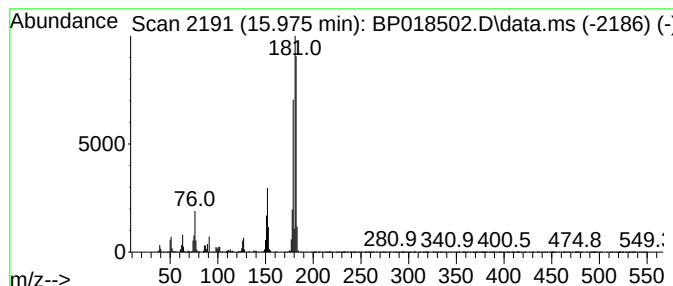
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 9H-Fluoren-3-ol Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.975	2.42 ng/ul	707759	Phenanthrene-d10	17.239

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			9H-Fluoren-3-ol	182	C13H10O	006344-67-8	60
2			2,4,6-Cycloheptatrien-1-one, 2-p...	182	C13H10O	014562-09-5	58
3			2,4,6-Cycloheptatrien-1-one, 2-p...	182	C13H10O	014562-09-5	58
4			9H-Fluoren-9-ol	182	C13H10O	001689-64-1	58
5			2-Hydroxyfluorene	182	C13H10O	002443-58-5	55



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP120123\
Data File : BP018502.D
Acq On : 02 Dec 2023 06:59
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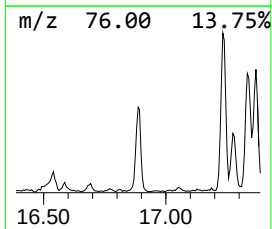
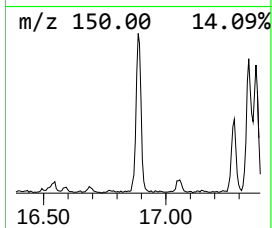
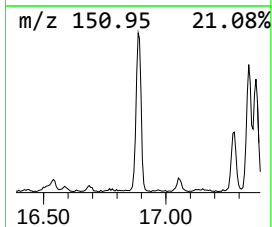
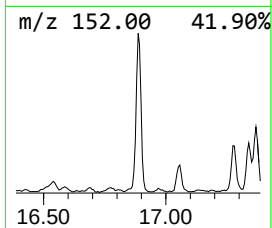
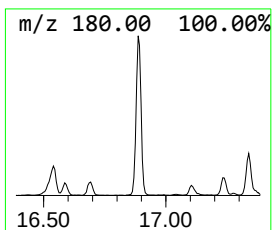
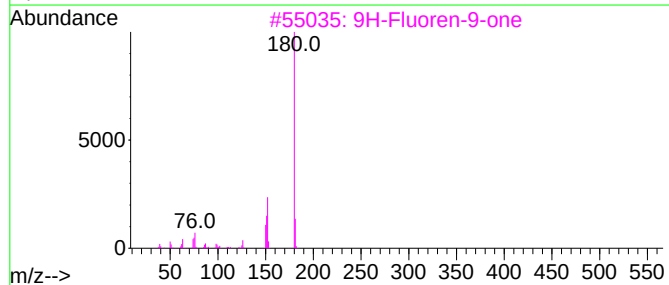
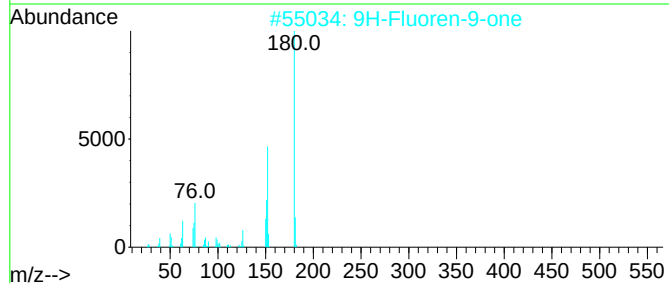
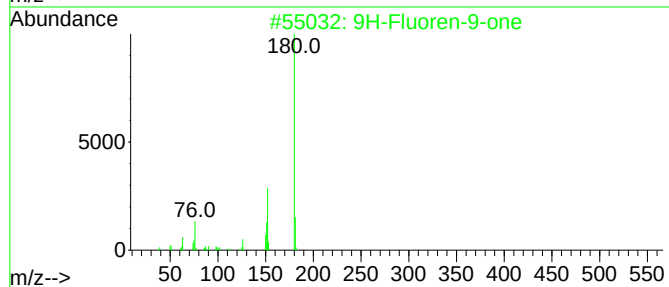
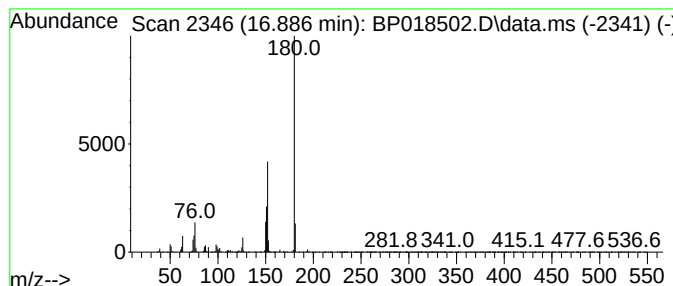
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 9H-Fluoren-9-one Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.886	2.07 ng/ul	605584	Phenanthrene-d10	17.239

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	9H-Fluoren-9-one			180	C13H8O	000486-25-9	96
2	9H-Fluoren-9-one			180	C13H8O	000486-25-9	95
3	9H-Fluoren-9-one			180	C13H8O	000486-25-9	95
4	9H-Fluoren-9-one			180	C13H8O	000486-25-9	94
5	9H-Fluoren-9-one			180	C13H8O	000486-25-9	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP120123\
Data File : BP018502.D
Acq On : 02 Dec 2023 06:59
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Misc :
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Instrument :
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ClientSampleId :
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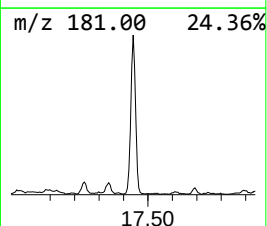
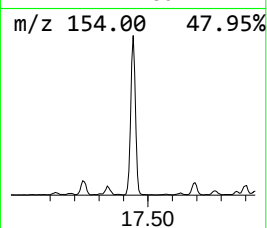
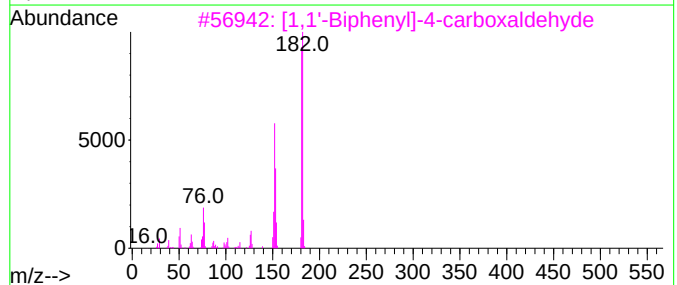
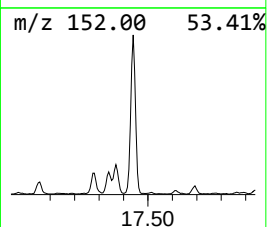
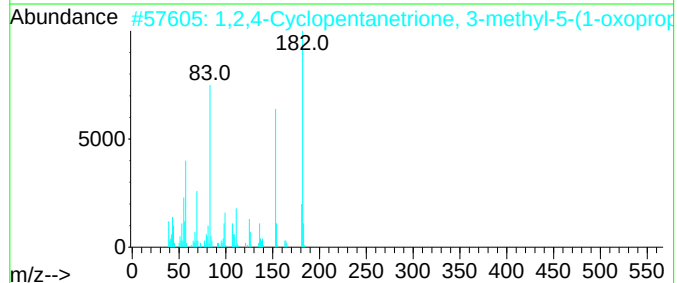
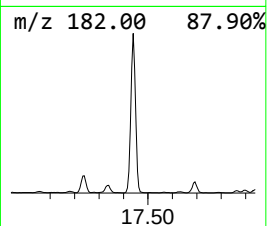
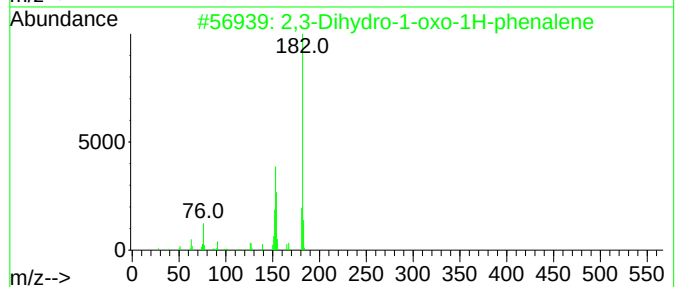
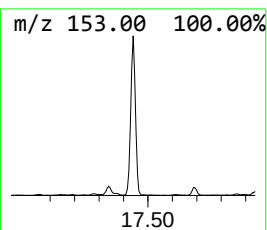
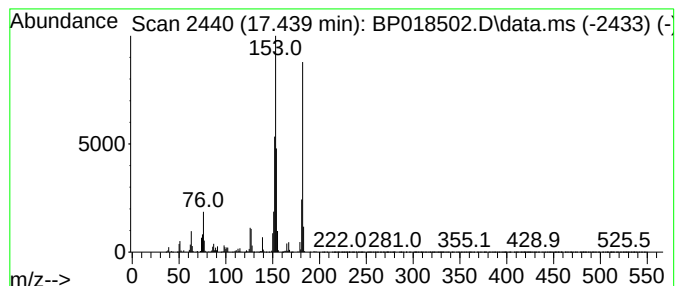
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 2,3-Dihydro-1-oxo-1H-phenalene Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.439	6.21 ng/ul	1815580	Phenanthrene-d10	17.239

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2,3-Dihydro-1-oxo-1H-phenalene	182	C13H10O	000518-85-4	76
2			1,2,4-Cyclopentanetrione, 3-meth...	182	C9H10O4	057174-14-8	50
3			[1,1'-Biphenyl]-4-carboxaldehyde	182	C13H10O	003218-36-8	49
4			9H-Fluoren-3-ol	182	C13H10O	006344-67-8	46
5			9H-Pyrido[3,4-b]indole, 1-methyl-	182	C12H10N2	000486-84-0	43



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP120123\
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Acq On : 02 Dec 2023 06:59
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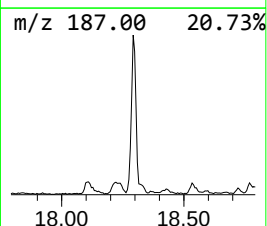
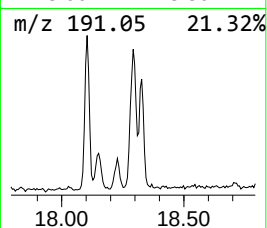
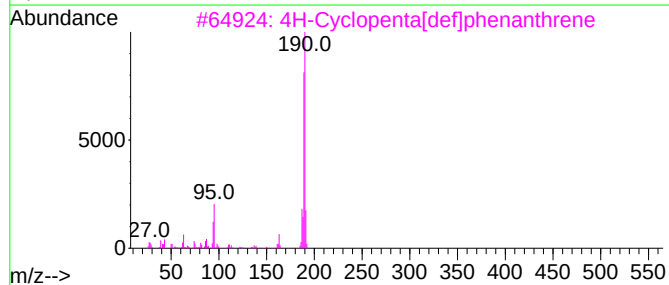
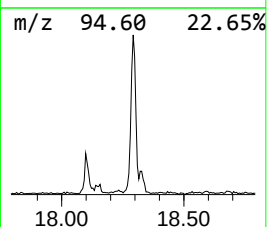
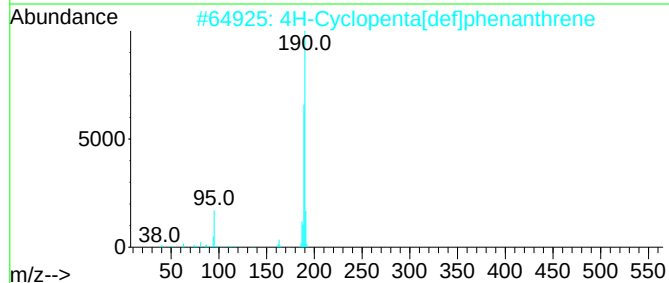
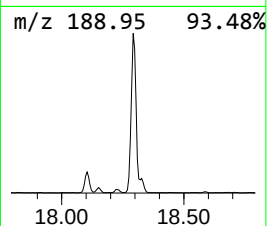
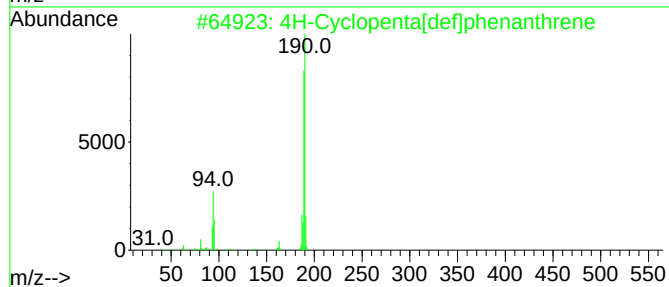
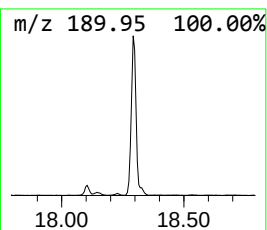
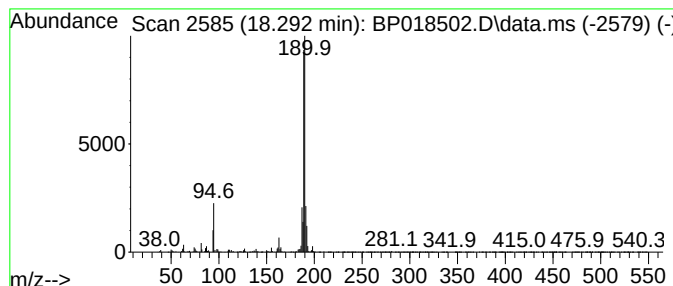
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 4H-Cyclopenta[def]phenanthrene Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.292	2.41 ng/ul	703765	Phenanthrene-d10	17.239

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	97
2			4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	87
3			4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	81
4			6H-Cyclobuta[jk]phenanthrene	190	C15H10	083469-43-6	56
5			Methyl diselenide	190	C2H6Se2	007101-31-7	49



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP120123\
Data File : BP018502.D
Acq On : 02 Dec 2023 06:59
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ALS Vial : 31 Sample Multiplier: 1

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ClientSampleId :
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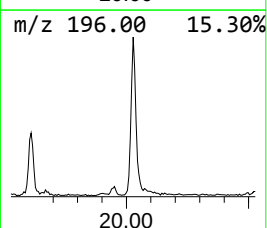
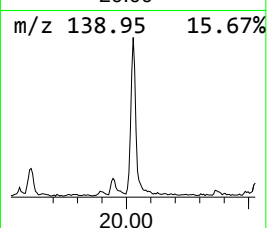
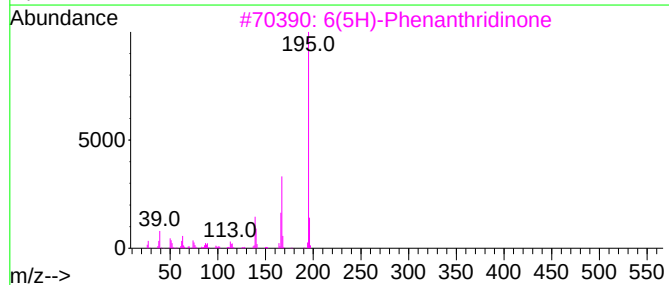
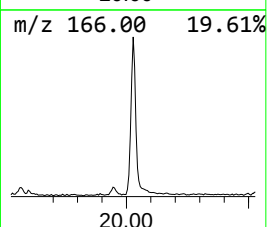
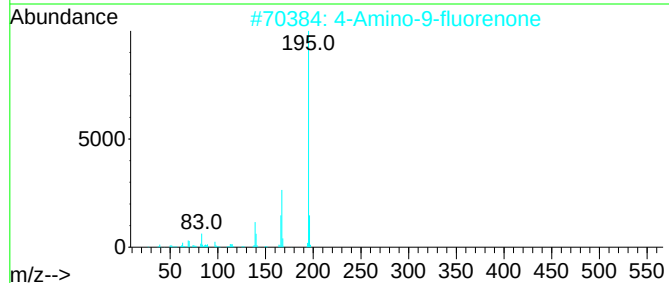
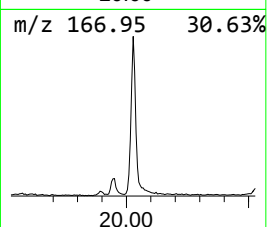
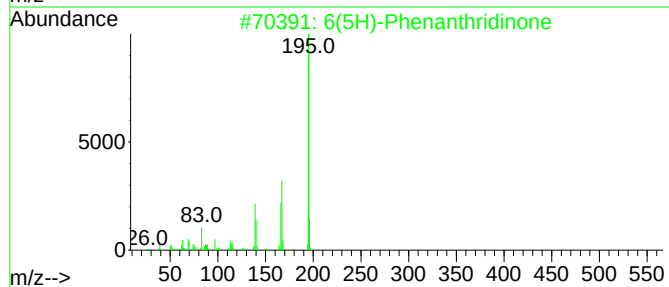
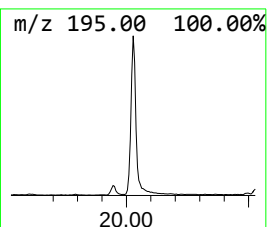
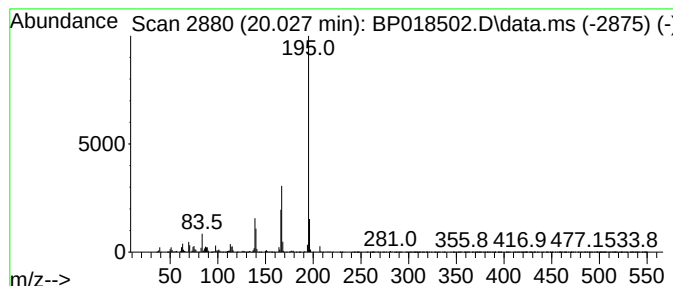
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 6(5H)-Phenanthridinone Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.027	2.59 ng/ul	713925	Chrysene-d12	21.322

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			6(5H)-Phenanthridinone	195	C13H9NO	001015-89-0	97
2			4-Amino-9-fluorenone	195	C13H9NO	004269-15-2	96
3			6(5H)-Phenanthridinone	195	C13H9NO	001015-89-0	96
4			9(10H)-Acridinone	195	C13H9NO	000578-95-0	95
5			9(10H)-Acridinone	195	C13H9NO	000578-95-0	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP120123\
Data File : BP018502.D
Acq On : 02 Dec 2023 06:59
Operator : MA/JU
Sample : 05332-16DL 5X
Misc :
ALS Vial : 31 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
YCG97DL

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120123.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
Benzene, 1-ethy...	8.969	2.0	ng/u1	258483	1	7.863	2546040	20.0
1H-Benzimidazol...	11.075	2.3	ng/u1	461599	2	10.657	4003260	20.0
3-Methylbenzoth...	12.504	2.0	ng/u1	406488	2	10.657	4003260	20.0
Naphthalene, 2,...	13.816	4.7	ng/u1	1264980	3	14.492	5342850	20.0
Naphthalene, 1,...	14.051	3.4	ng/u1	906401	3	14.492	5342850	20.0
9H-Fluoren-3-ol	15.975	2.4	ng/u1	707759	4	17.239	5847940	20.0
9H-Fluoren-9-one	16.886	2.1	ng/u1	605584	4	17.239	5847940	20.0
2,3-Dihydro-1-o...	17.439	6.2	ng/u1	1815580	4	17.239	5847940	20.0
4H-Cyclopenta[d...	18.292	2.4	ng/u1	703765	4	17.239	5847940	20.0
6(5H)-Phenanthr...	20.027	2.6	ng/u1	713925	5	21.322	5523410	20.0