

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP012324\
 Data File : BP019430.D
 Acq On : 23 Jan 2024 21:16
 Operator : MA/JU
 Sample : P1158-10
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0AK6

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-BP010824.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BP019430.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.128	3	7	12	rVB	877649	885594	24.26%	2.694%
2	3.464	61	64	69	rVB	20459	25913	0.71%	0.079%
3	3.946	143	146	153	rVB2	16227	22850	0.63%	0.070%
4	5.269	367	371	379	rVB	29363	42399	1.16%	0.129%
5	5.405	388	394	406	rVB2	11898	24975	0.68%	0.076%
6	5.775	452	457	461	rVB	11576	19357	0.53%	0.059%
7	6.246	531	537	543	rBV5	7624	14827	0.41%	0.045%
8	6.746	618	622	629	rVB7	6437	10208	0.28%	0.031%
9	6.852	635	640	644	rBV2	20610	32723	0.90%	0.100%
10	6.910	644	650	655	rVV	20782	38816	1.06%	0.118%
11	6.987	658	663	669	rVB2	10103	16346	0.45%	0.050%
12	7.152	684	691	699	rBV6	11595	31595	0.87%	0.096%
13	7.228	699	704	706	rVV6	6087	10328	0.28%	0.031%
14	7.316	713	719	725	rVV6	5778	12463	0.34%	0.038%
15	7.410	725	735	741	rVB3	20639	42240	1.16%	0.128%
16	7.657	769	777	785	rBV3	8531	24983	0.68%	0.076%
17	7.769	785	796	809	rVV	660210	1077631	29.52%	3.278%
18	7.893	809	817	822	rVV4	19549	47512	1.30%	0.145%
19	8.134	851	858	865	rVB	25563	44258	1.21%	0.135%
20	8.305	876	887	892	rBV	7895	24305	0.67%	0.074%
21	8.405	899	904	910	rBV3	14359	27793	0.76%	0.085%
22	8.834	971	977	979	rVV7	5301	9139	0.25%	0.028%
23	8.863	979	982	988	rVV3	14063	24538	0.67%	0.075%
24	8.987	998	1003	1012	rVB	17859	32538	0.89%	0.099%
25	9.110	1021	1024	1030	rVB8	7566	12834	0.35%	0.039%
26	9.393	1066	1072	1076	rBV5	7284	14006	0.38%	0.043%
27	9.452	1078	1082	1089	rVB5	9802	19367	0.53%	0.059%
28	9.810	1137	1143	1151	rVB7	7389	16798	0.46%	0.051%
29	9.957	1163	1168	1175	rBV4	15862	30871	0.85%	0.094%
30	10.422	1242	1247	1254	rBV4	7105	17803	0.49%	0.054%
31	10.563	1260	1271	1275	rBV	868712	1479955	40.55%	4.502%
32	10.610	1275	1279	1286	rVV	315408	517303	14.17%	1.574%
33	10.675	1286	1290	1292	rVV5	7079	10787	0.30%	0.033%
34	10.722	1294	1298	1305	rVV6	8868	17211	0.47%	0.052%
35	11.475	1421	1426	1431	rVB9	8762	16883	0.46%	0.051%
36	11.581	1439	1444	1449	rBV3	13354	25403	0.70%	0.077%

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Integration Parameters: LSCINT.P

Integrator: RTE

Smoothing : OFF

Filtering: 5

Sampling : 1

Min Area: 0.1 % of largest Peak

Start Thrs: 0.2

Max Peaks: 100

Stop Thrs : 0

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

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Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP010824.MA.M
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37	11.981	1508	1512	1516	rBV6	11877	19293	0.53%	0.059%
38	12.228	1547	1554	1561	rVB	124870	207767	5.69%	0.632%
39	12.445	1583	1591	1601	rBV	105963	188447	5.16%	0.573%
40	12.610	1615	1619	1623	rBV7	6512	11005	0.30%	0.033%

41	12.834	1654	1657	1662	rVB6	12999	18306	0.50%	0.056%
42	12.893	1663	1667	1669	rBV5	6430	10807	0.30%	0.033%
43	12.940	1670	1675	1680	rVV3	23249	36827	1.01%	0.112%
44	12.993	1680	1684	1690	rVB9	9980	18297	0.50%	0.056%
45	13.175	1710	1715	1719	rBV5	13883	29037	0.80%	0.088%

46	13.245	1721	1727	1732	rVB3	39800	78019	2.14%	0.237%
47	13.440	1757	1760	1763	rBV	18099	26329	0.72%	0.080%
48	13.569	1778	1782	1784	rBV	35247	47542	1.30%	0.145%
49	13.734	1804	1810	1815	rBV2	77149	152469	4.18%	0.464%
50	13.787	1815	1819	1822	rVV	52893	89070	2.44%	0.271%

51	13.816	1822	1824	1829	rVV	51248	61401	1.68%	0.187%
52	13.887	1832	1836	1840	rVV3	54034	80679	2.21%	0.245%
53	13.934	1841	1844	1846	rVV4	21545	32505	0.89%	0.099%
54	13.969	1847	1850	1853	rVV2	40951	53401	1.46%	0.162%
55	14.004	1853	1856	1860	rVB5	20336	27126	0.74%	0.083%

56	14.134	1874	1878	1885	rVV3	49192	85394	2.34%	0.260%
57	14.228	1890	1894	1902	rVB4	53646	91233	2.50%	0.278%
58	14.310	1902	1908	1912	rBV2	36443	61586	1.69%	0.187%
59	14.410	1919	1925	1933	rBV2	1326946	2116267	57.98%	6.438%
60	14.475	1933	1936	1944	rVB	111004	172831	4.73%	0.526%

61	14.634	1957	1963	1970	rVB4	27319	72897	2.00%	0.222%
62	14.728	1972	1979	1982	rBV7	21588	52172	1.43%	0.159%
63	14.816	1989	1994	1999	rVV	186957	265220	7.27%	0.807%
64	14.887	2003	2006	2010	rBV	29215	43325	1.19%	0.132%
65	14.928	2011	2013	2018	rVB5	27763	35975	0.99%	0.109%

66	15.034	2027	2031	2037	rBV5	44695	96207	2.64%	0.293%
67	15.192	2053	2058	2062	rBV6	47997	87832	2.41%	0.267%
68	15.463	2098	2104	2109	rBV	413965	613673	16.81%	1.867%
69	15.628	2128	2132	2135	rBV4	29277	43990	1.21%	0.134%
70	15.669	2135	2139	2144	rVB3	80032	120822	3.31%	0.368%

71	15.757	2151	2154	2161	rVB2	85199	137950	3.78%	0.420%
72	15.875	2172	2174	2177	rBV3	22289	31616	0.87%	0.096%
73	15.910	2177	2180	2187	rVB3	46377	91049	2.49%	0.277%
74	16.145	2213	2220	2225	rBV2	197761	363731	9.96%	1.106%
75	16.469	2270	2275	2279	rBV3	64956	109334	3.00%	0.333%

76	16.981	2357	2362	2369	rVB2	213358	402830	11.04%	1.225%
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77	17.045	2370	2373	2377	rBV3	75150	96863	2.65%	0.295%
78	17.169	2388	2394	2397	rBV	1619843	2219602	60.81%	6.752%
79	17.210	2397	2401	2408	rVB	2567493	3650150	100.00%	11.104%
80	17.304	2414	2417	2420	rBV	58347	69552	1.91%	0.212%
81	17.586	2461	2465	2469	rBV2	80482	144525	3.96%	0.440%
82	17.757	2491	2494	2502	rVB2	168279	345725	9.47%	1.052%
83	18.039	2539	2542	2546	rBV	170112	250004	6.85%	0.761%
84	18.086	2547	2550	2556	rVB	278871	394641	10.81%	1.201%
85	18.228	2570	2574	2578	rBV3	514719	737513	20.21%	2.244%
86	18.504	2618	2621	2622	rBV2	94433	103611	2.84%	0.315%
87	18.528	2623	2625	2629	rVB	170135	163679	4.48%	0.498%
88	18.757	2661	2664	2667	rBV2	135470	161644	4.43%	0.492%
89	18.933	2691	2694	2698	rBV	112056	169377	4.64%	0.515%
90	19.027	2707	2710	2712	rBV2	134012	154589	4.24%	0.470%
91	19.186	2733	2737	2742	rBV	2715825	3587624	98.29%	10.914%
92	19.239	2742	2746	2751	rVV	755368	901980	24.71%	2.744%
93	19.327	2757	2761	2767	rBV5	182058	248303	6.80%	0.755%
94	19.516	2789	2793	2794	rBV2	332723	424074	11.62%	1.290%
95	19.539	2794	2797	2805	rVB	1347040	1847028	50.60%	5.619%
96	19.945	2863	2866	2870	rBV	190469	223761	6.13%	0.681%
97	20.122	2894	2896	2900	rVB	251074	284042	7.78%	0.864%
98	21.269	3086	3091	3095	rBV2	1568857	2844650	77.93%	8.654%
99	21.310	3096	3098	3106	rVB	891497	997338	27.32%	3.034%
100	23.621	3488	3491	3498	rVB	1047362	1845313	50.55%	5.614%

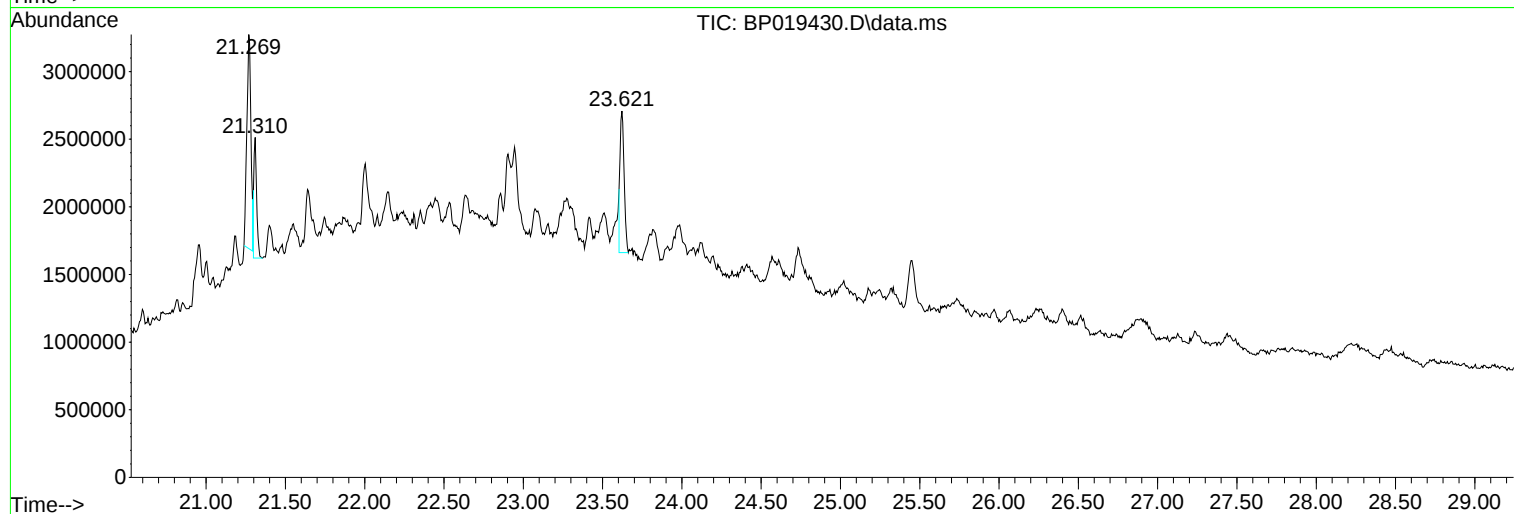
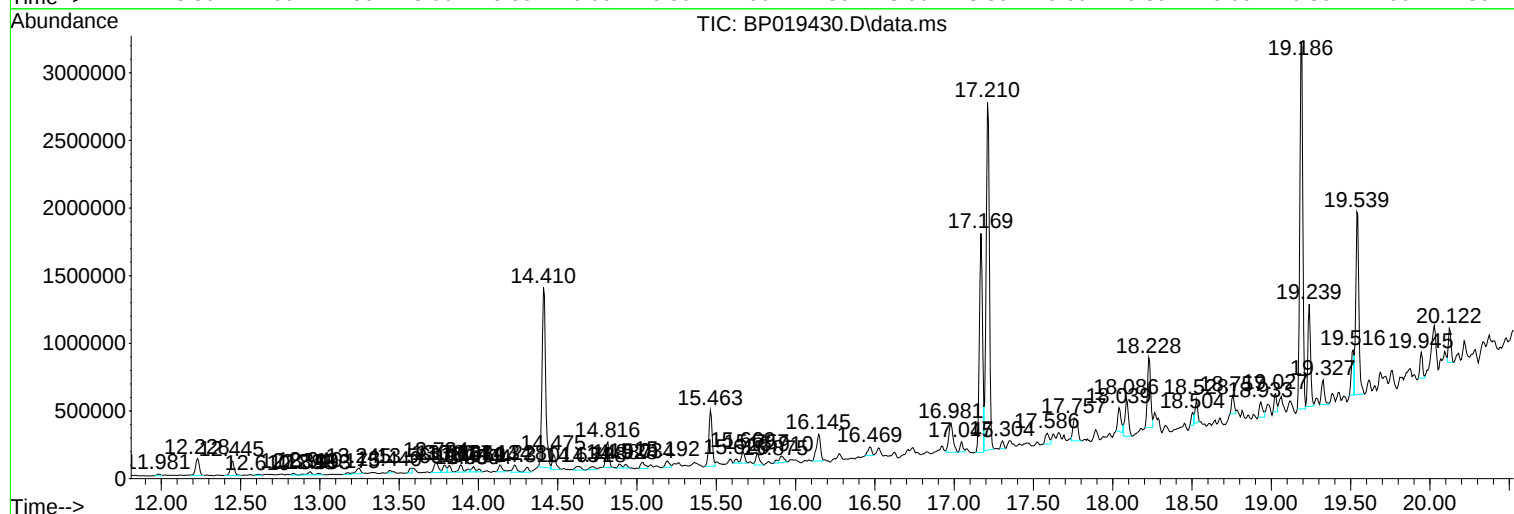
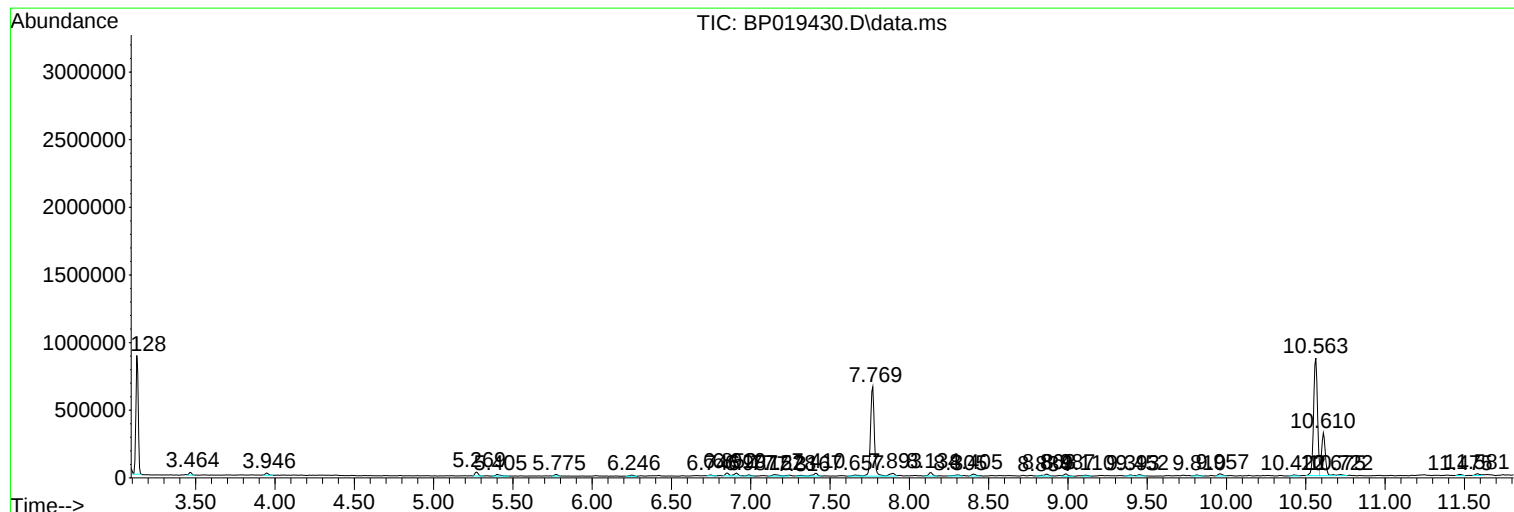
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Instrument :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP010824.MA.M
Quant Title : SVOA CALIBRATION

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



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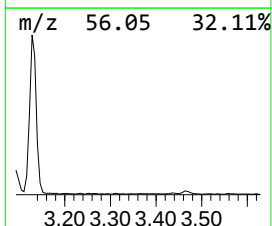
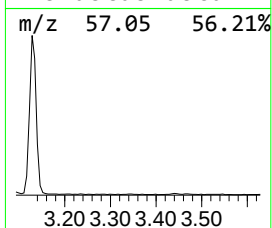
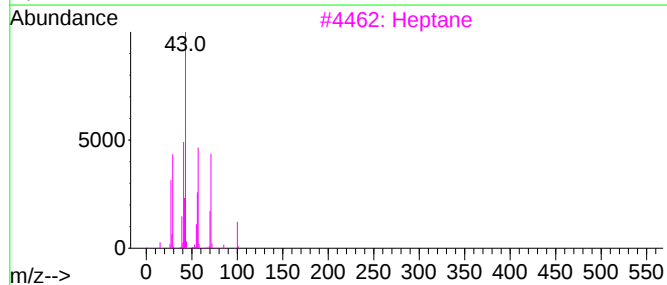
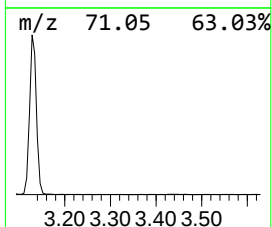
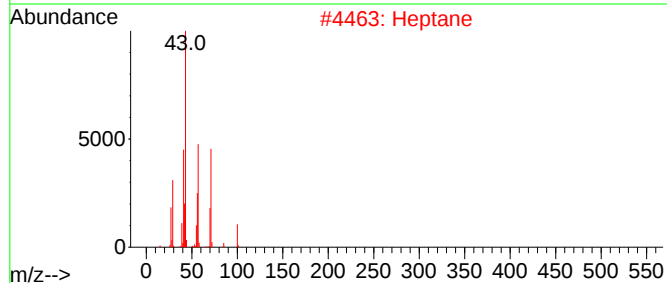
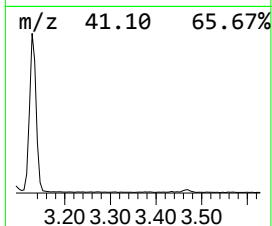
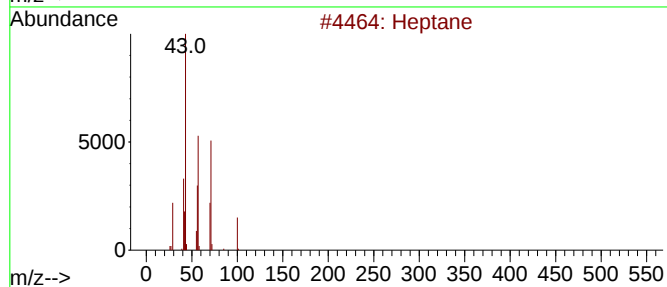
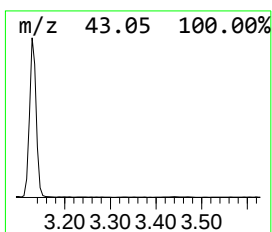
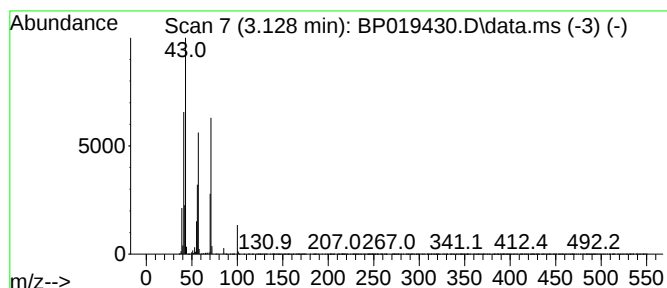
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP010824.MA.M
Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.128	16.44 ng/ul	885594	1,4-Dichlorobenzene-d4	7.769

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptane			100	C7H16	000142-82-5	91
2	Heptane			100	C7H16	000142-82-5	90
3	Heptane			100	C7H16	000142-82-5	90
4	Heptane			100	C7H16	000142-82-5	86
5	Pentane, 2-methyl-			86	C6H14	000107-83-5	47



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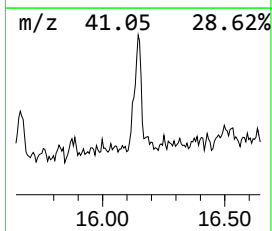
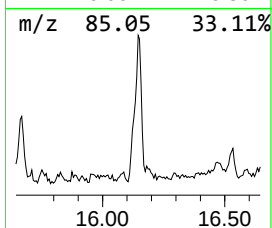
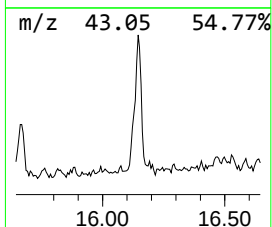
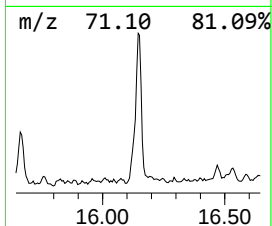
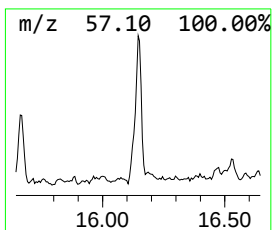
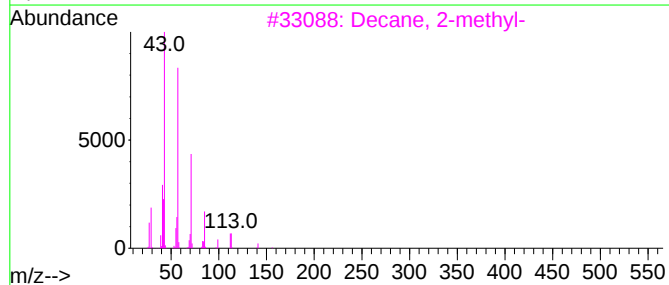
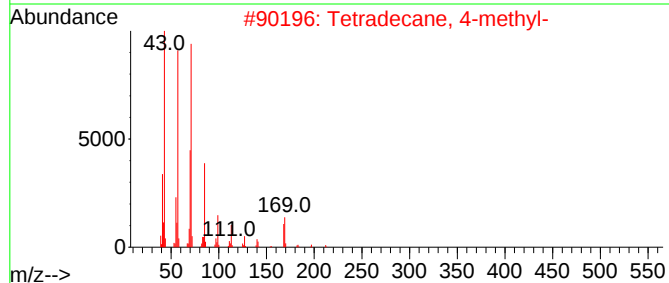
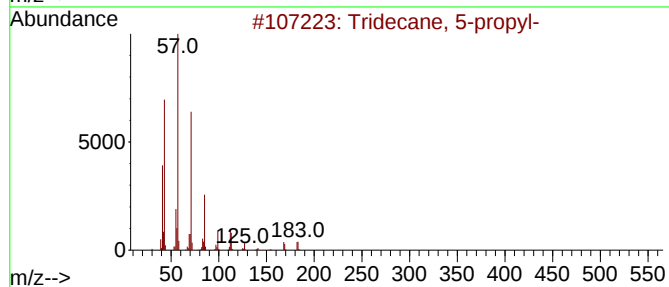
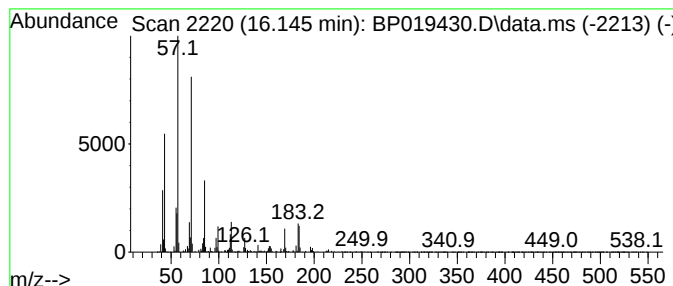
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Quant Title : SVOA CALIBRATION

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

Peak Number 2 (DEL) Alkane: Straight-Chai... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.	
16.145	3.28 ng/ul	363731	Phenanthrene-d10	17.169	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tridecane, 5-propyl-	226	C16H34	055045-11-9	91
2	Tetradecane, 4-methyl-	212	C15H32	025117-24-2	80
3	Decane, 2-methyl-	156	C11H24	006975-98-0	72
4	Heptacosane	380	C27H56	000593-49-7	72
5	2-Bromo dodecane	248	C12H25Br	013187-99-0	64



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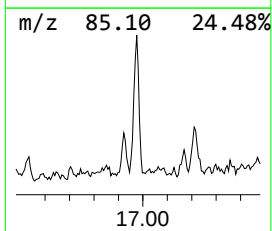
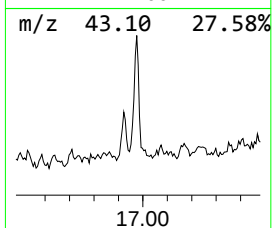
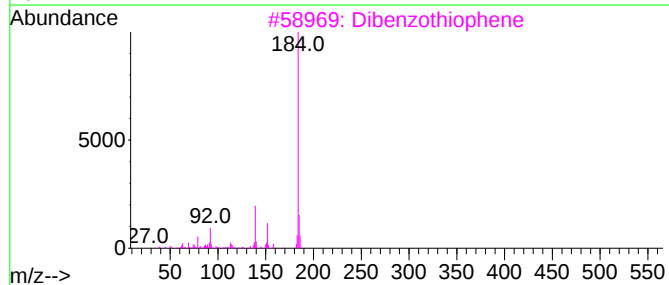
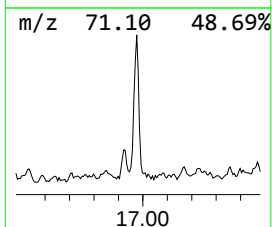
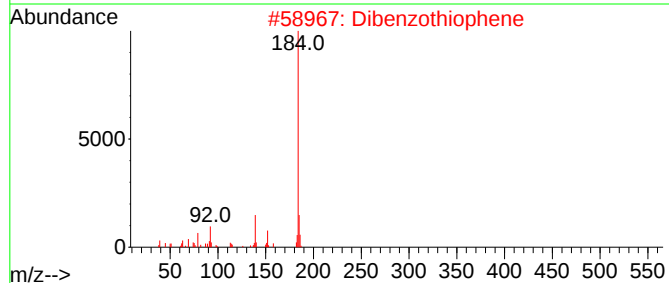
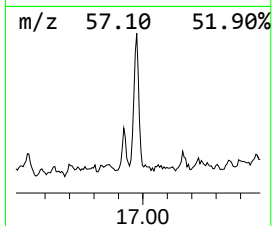
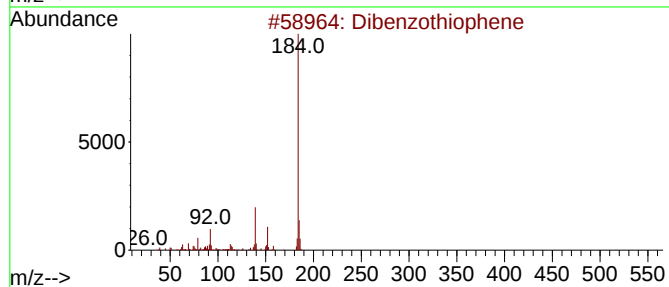
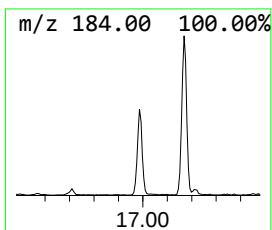
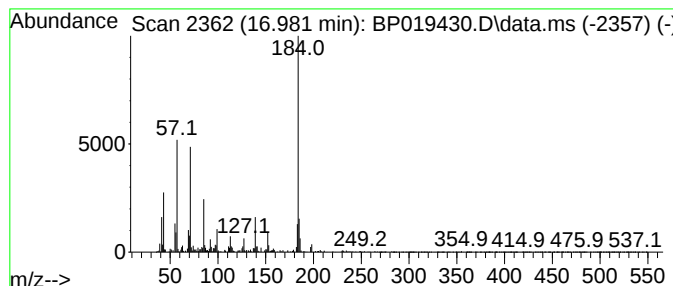
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP010824.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 Dibenzothiophene Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.980	3.63 ng/ul	402830	Phenanthrene-d10	17.169

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Dibenzothiophene	184	C12H8S	000132-65-0	60
2			Dibenzothiophene	184	C12H8S	000132-65-0	60
3			Dibenzothiophene	184	C12H8S	000132-65-0	60
4			Dibenzothiophene	184	C12H8S	000132-65-0	55
5			Dibenzothiophene	184	C12H8S	000132-65-0	55



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP012324\
Data File : BP019430.D
Acq On : 23 Jan 2024 21:16
Operator : MA/JU
Sample : P1158-10
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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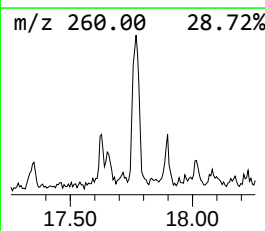
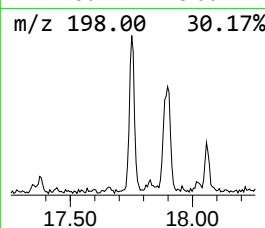
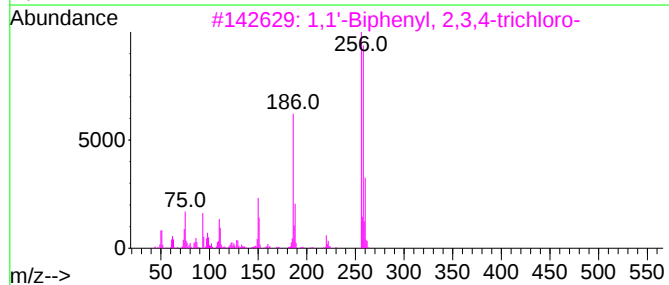
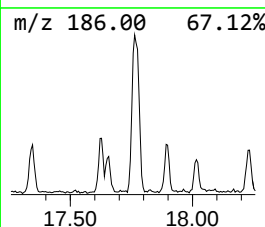
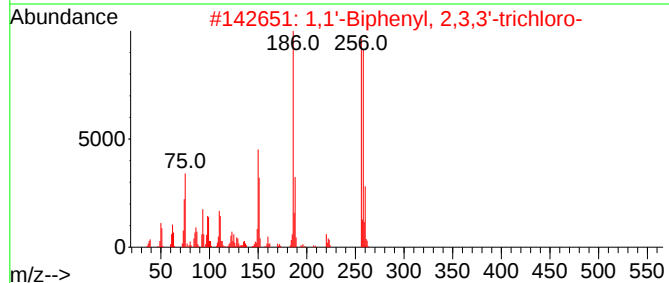
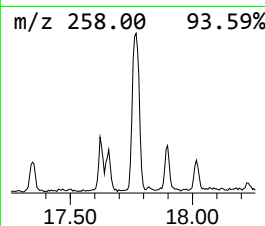
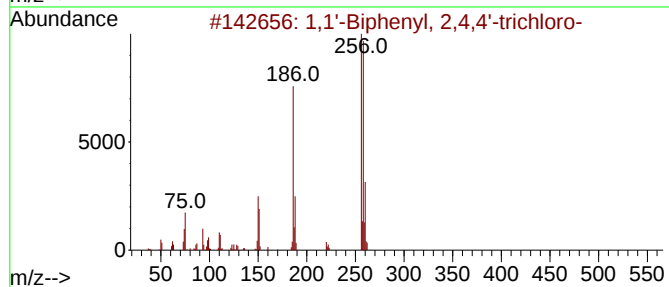
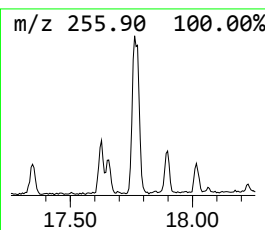
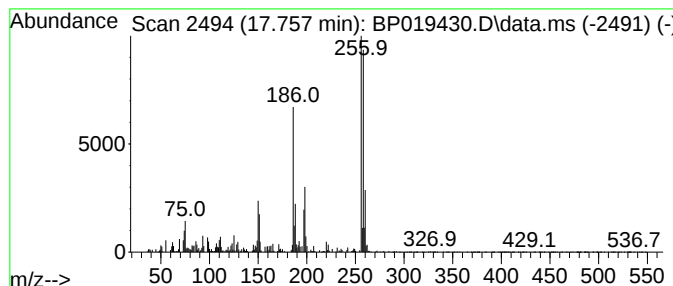
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 1,1'-Biphenyl, 2,4,4'-trich... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.757	3.12 ng/ul	345725	Phenanthrene-d10	17.169

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl, 2,4,4'-trichloro-	256	C12H7Cl3	007012-37-5	99		
2	1,1'-Biphenyl, 2,3,3'-trichloro-	256	C12H7Cl3	038444-84-7	99		
3	1,1'-Biphenyl, 2,3,4-trichloro-	256	C12H7Cl3	055702-46-0	95		
4	2,4',6-Trichloro-1,1'-biphenyl	256	C12H7Cl3	038444-77-8	94		
5	1,1'-Biphenyl, 2,4,6-trichloro-	256	C12H7Cl3	035693-92-6	94		



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP012324\
Data File : BP019430.D
Acq On : 23 Jan 2024 21:16
Operator : MA/JU
Sample : P1158-10
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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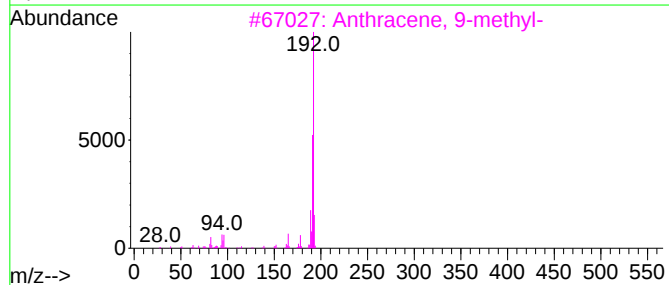
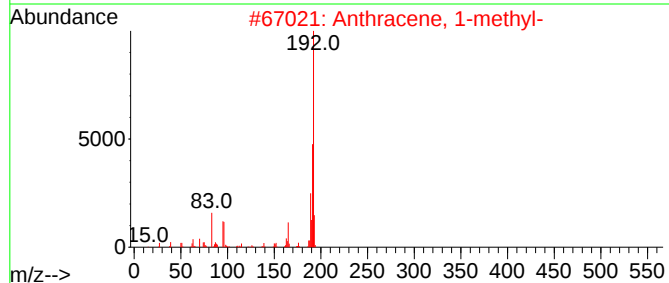
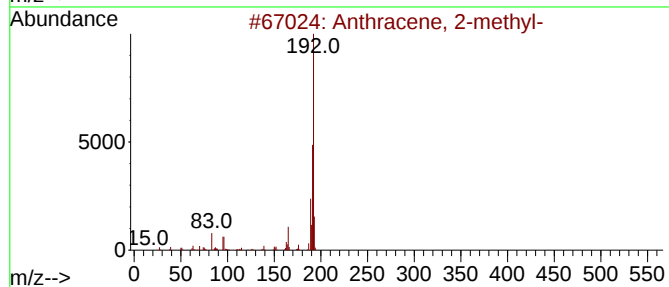
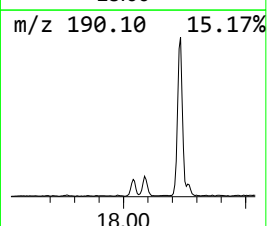
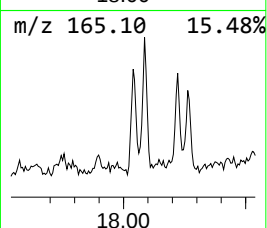
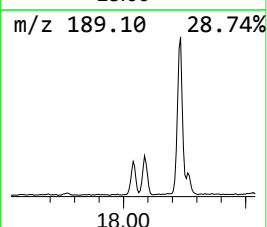
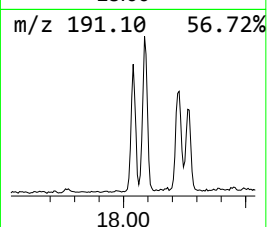
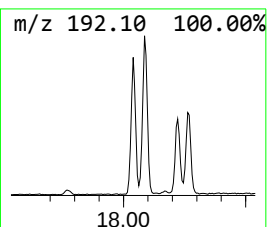
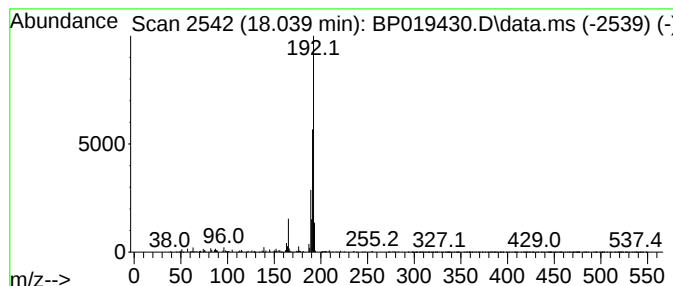
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 Anthracene, 2-methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.039	2.25 ng/ul	250004	Phenanthrene-d10	17.169

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Anthracene, 2-methyl-	192	C15H12	000613-12-7	87
2			Anthracene, 1-methyl-	192	C15H12	000610-48-0	87
3			Anthracene, 9-methyl-	192	C15H12	000779-02-2	87
4			1H-Indene, 2-phenyl-	192	C15H12	004505-48-0	87
5			Anthracene, 2-methyl-	192	C15H12	000613-12-7	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP012324\
Data File : BP019430.D
Acq On : 23 Jan 2024 21:16
Operator : MA/JU
Sample : P1158-10
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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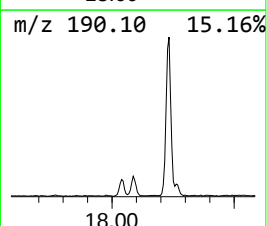
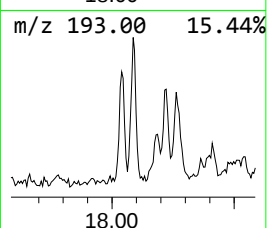
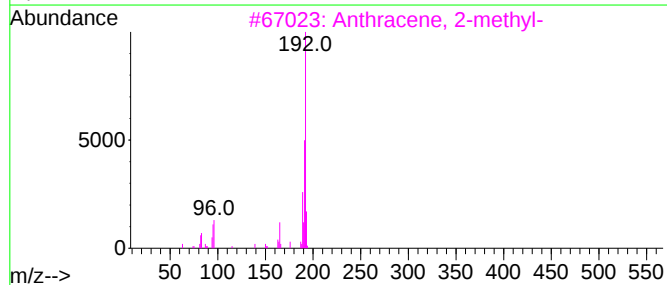
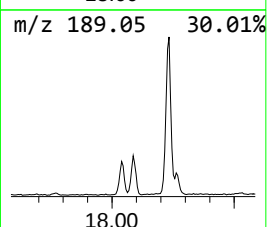
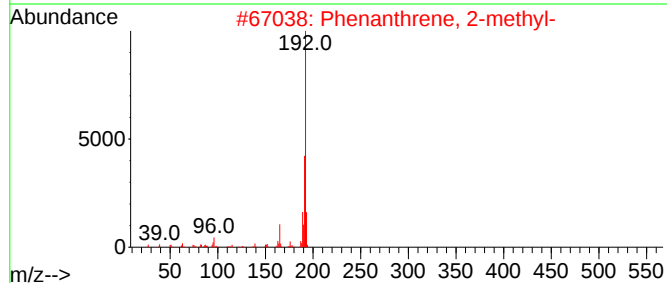
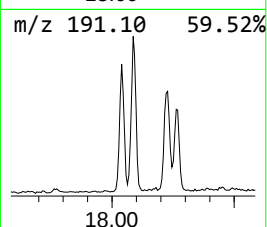
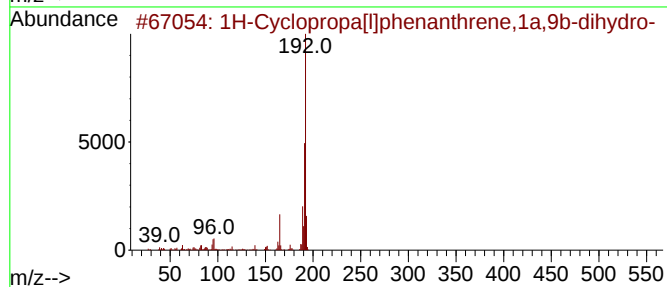
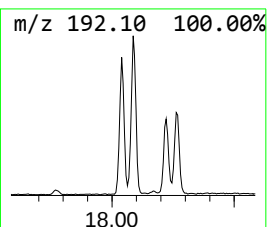
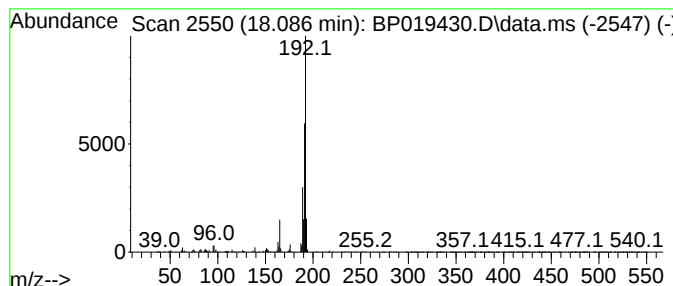
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 1H-Cyclopropa[l]phenanthren... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.086	3.56 ng/ul	394641	Phenanthrene-d10	17.169

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1H-Cyclopropa[l]phenanthrene,1a,...	192	C15H12	000949-41-7	93
2			Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	93
3			Anthracene, 2-methyl-	192	C15H12	000613-12-7	93
4			Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	93
5			Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP012324\
Data File : BP019430.D
Acq On : 23 Jan 2024 21:16
Operator : MA/JU
Sample : P1158-10
Misc :
ALS Vial : 17 Sample Multiplier: 1

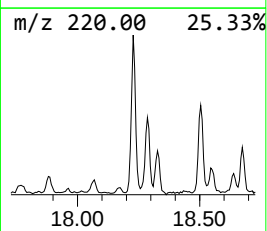
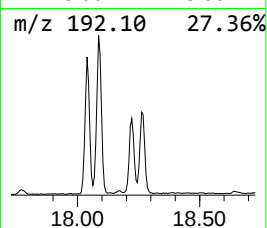
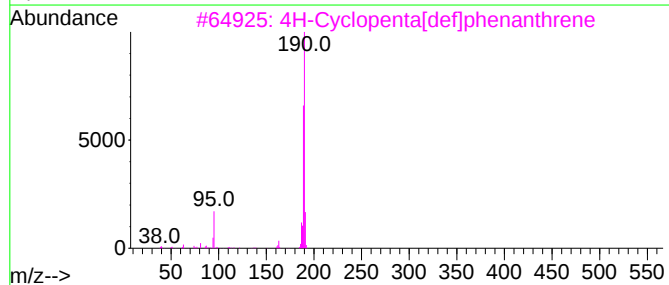
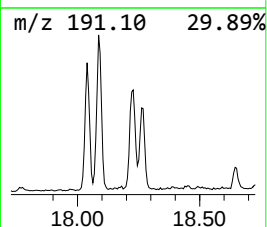
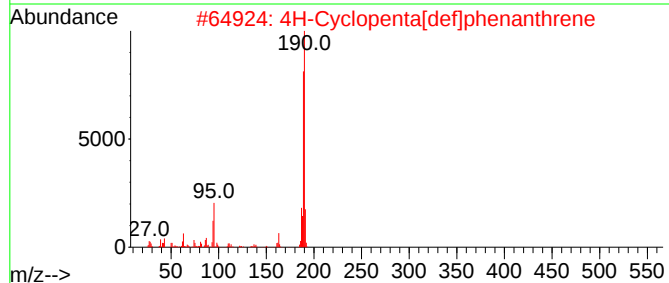
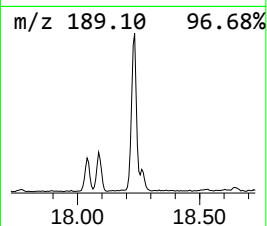
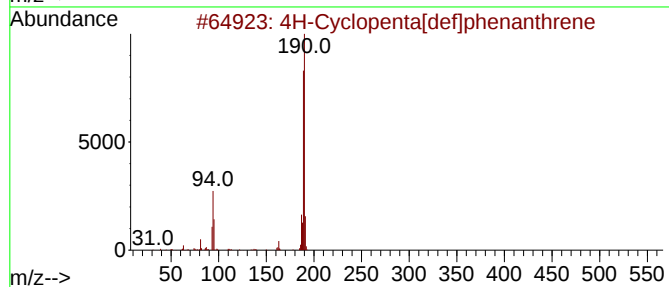
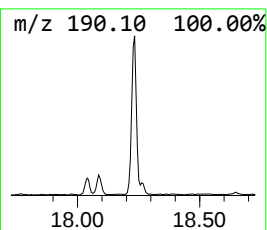
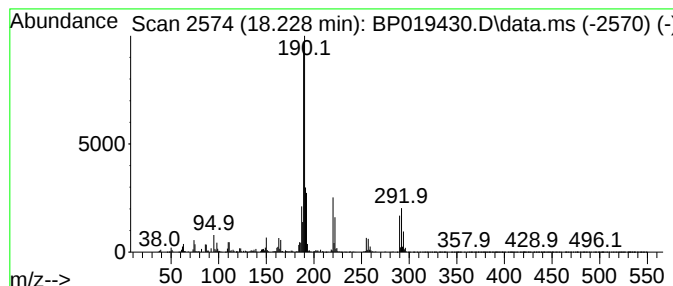
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ClientSampleId :
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 4H-Cyclopenta[def]phenanthrene Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.	
18.227	6.65 ng/ul	737513	Phenanthrene-d10	17.169	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	58
2	4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	50
3	4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	50
4	2,2'-Bis(4,5-dimethylimidazole)	190	C10H14N4	069286-06-2	47
5	4-(4-Methyl-piperidin-1-yl)-phen...	190	C12H18N2	342013-25-6	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP012324\
Data File : BP019430.D
Acq On : 23 Jan 2024 21:16
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Sample : P1158-10
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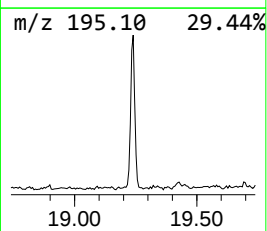
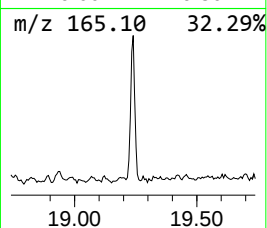
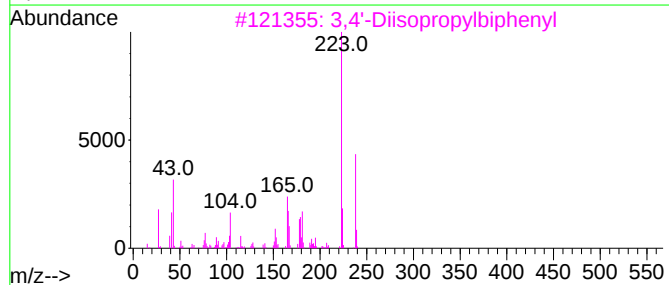
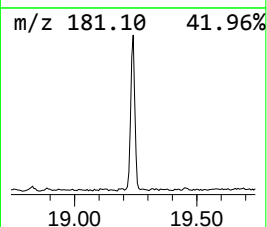
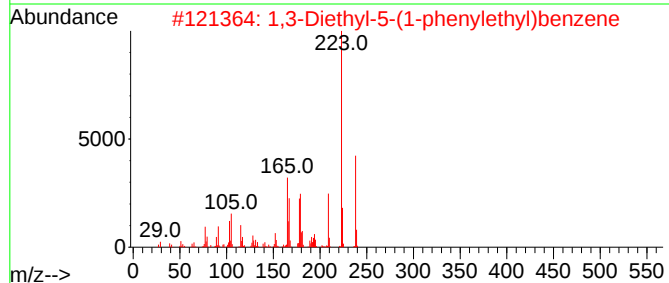
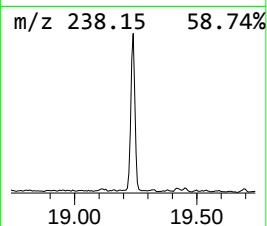
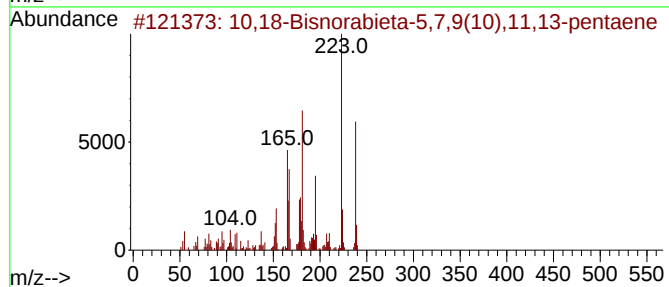
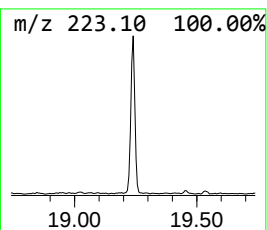
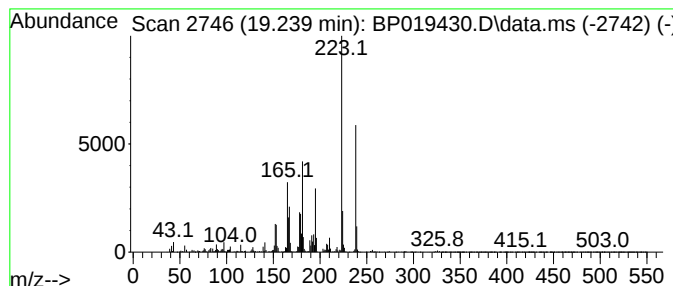
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 10,18-Bisnorabieta-5,7,9(10... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.239	6.34 ng/ul	901980	Chrysene-d12	21.274

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	10,18-Bisnorabieta-5,7,9(10),11,...	238	C18H22	006566-19-4	99		
2	1,3-Diethyl-5-(1-phenylethyl)ben...	238	C18H22	1000482-32-6	86		
3	3,4'-Diisopropylbiphenyl	238	C18H22	061434-46-6	86		
4	12-Azabicyclo[9.2.2]pentadecan-1...	223	C14H25NO	1000191-26-7	74		
5	cyclopropanone, 2-(4-methoxyphen...	238	C16H14O2	1000401-16-2	62		



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP012324\
Data File : BP019430.D
Acq On : 23 Jan 2024 21:16
Operator : MA/JU
Sample : P1158-10
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP010824.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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(DEL) Alkane: S...	16.145	3.3	ng/ul	363731	4	17.169	2219600	20.0
Dibenzothiophene	16.980	3.6	ng/ul	402830	4	17.169	2219600	20.0
1,1'-Biphenyl, ...	17.757	3.1	ng/ul	345725	4	17.169	2219600	20.0
Anthracene, 2-m...	18.039	2.3	ng/ul	250004	4	17.169	2219600	20.0
1H-Cyclopropa[l...	18.086	3.6	ng/ul	394641	4	17.169	2219600	20.0
4H-Cyclopenta[d...	18.227	6.7	ng/ul	737513	4	17.169	2219600	20.0
10,18-Bisnorabi...	19.239	6.3	ng/ul	901980	5	21.274	2844650	20.0