

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP080624\
 Data File : BP021447.D
 Acq On : 08 Aug 2024 12:38
 Operator : CG/JU
 Sample : P3437-14DL 10X
 Misc :
 ALS Vial : 39 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 E05Y4DL

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-BP071024.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BP021447.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.669	114	116	123	rVB5	9618	16317	0.26%	0.034%
2	3.875	148	151	154	rVB4	5581	6502	0.10%	0.014%
3	7.040	677	689	691	rBV4	14035	35625	0.57%	0.075%
4	7.281	722	730	743	rBV4	14364	52746	0.84%	0.111%
5	7.622	781	788	812	rBV	251494	778454	12.40%	1.640%
6	8.181	877	883	889	rBV8	4057	8491	0.14%	0.018%
7	8.487	927	935	936	rBV4	10121	20228	0.32%	0.043%
8	8.863	991	999	1004	rBV4	16439	44089	0.70%	0.093%
9	9.551	1108	1116	1117	rBV6	9086	17198	0.27%	0.036%
10	10.192	1216	1225	1233	rBV6	15406	64312	1.02%	0.135%
11	10.381	1249	1257	1260	rBV	485940	982509	15.66%	2.070%
12	10.434	1260	1266	1283	rVV	654938	1717459	27.37%	3.618%
13	10.563	1285	1288	1297	rVB3	15649	33128	0.53%	0.070%
14	10.645	1297	1302	1305	rBV5	7519	15006	0.24%	0.032%
15	11.245	1402	1404	1419	rVB5	5189	15362	0.24%	0.032%
16	11.957	1520	1525	1530	rBV3	6072	11510	0.18%	0.024%
17	12.051	1534	1541	1561	rBV	395405	881329	14.04%	1.856%
18	12.281	1570	1580	1597	rVB	175575	452023	7.20%	0.952%
19	13.045	1704	1710	1712	rBV7	4929	7415	0.12%	0.016%
20	13.086	1712	1717	1732	rVV	131261	265701	4.23%	0.560%
21	13.281	1744	1750	1764	rVB	43420	104851	1.67%	0.221%
22	13.434	1774	1776	1792	rVB2	67457	119017	1.90%	0.251%
23	13.569	1792	1799	1804	rBV2	86808	177592	2.83%	0.374%
24	13.628	1804	1809	1813	rVV2	59851	111668	1.78%	0.235%
25	13.681	1813	1818	1831	rVV	74042	180498	2.88%	0.380%
26	13.816	1835	1841	1857	rVB4	36089	104182	1.66%	0.219%
27	13.951	1857	1864	1882	rBV2	113825	259650	4.14%	0.547%
28	14.251	1908	1915	1921	rBV	1006855	1651656	26.32%	3.479%
29	14.322	1921	1927	1945	rVB	945458	1682893	26.82%	3.545%
30	14.481	1946	1954	1963	rBV3	32808	90323	1.44%	0.190%
31	14.575	1965	1970	1976	rVB3	29983	50867	0.81%	0.107%
32	14.669	1978	1986	1996	rBV	572217	1124454	17.92%	2.369%
33	14.745	1996	1999	2004	rVB2	24192	34867	0.56%	0.073%
34	14.892	2018	2024	2030	rVV3	16191	30518	0.49%	0.064%
35	14.963	2031	2036	2046	rVV4	17974	32578	0.52%	0.069%
36	15.063	2046	2053	2057	rVV3	16854	35816	0.57%	0.075%

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Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
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37	15.104	2057	2060	2067	rVB5	16225	26778	0.43%	0.056%
38	15.222	2076	2080	2084	rBV5	6740	12534	0.20%	0.026%
39	15.280	2084	2090	2094	rVV	142540	244628	3.90%	0.515%
40	15.333	2094	2099	2112	rVV	707937	1358473	21.65%	2.862%

41	15.457	2112	2120	2123	rVV4	50155	140270	2.24%	0.295%
42	15.492	2123	2126	2131	rVV3	98223	170829	2.72%	0.360%
43	15.545	2131	2135	2138	rVV5	30520	58869	0.94%	0.124%
44	15.592	2138	2143	2147	rVV3	43341	91809	1.46%	0.193%
45	15.651	2147	2153	2164	rVB	95537	180848	2.88%	0.381%

46	15.798	2172	2178	2189	rBV	83245	235475	3.75%	0.496%
47	15.892	2189	2194	2202	rVB4	22741	50359	0.80%	0.106%
48	15.992	2206	2211	2223	rBV7	18354	37715	0.60%	0.079%
49	16.151	2231	2238	2245	rBV	42239	68305	1.09%	0.144%
50	16.322	2263	2267	2268	rBV4	9524	11552	0.18%	0.024%

51	16.375	2268	2276	2280	rVV	73033	131756	2.10%	0.278%
52	16.422	2280	2284	2288	rVV2	30655	46310	0.74%	0.098%
53	16.463	2288	2291	2294	rVV5	6840	10637	0.17%	0.022%
54	16.516	2296	2300	2304	rVB	27747	40922	0.65%	0.086%
55	16.563	2304	2308	2310	rBV5	12469	15357	0.24%	0.032%

56	16.592	2310	2313	2316	rVV2	23177	29432	0.47%	0.062%
57	16.633	2316	2320	2323	rVB2	22399	29863	0.48%	0.063%
58	16.810	2343	2350	2356	rBV5	39136	77652	1.24%	0.164%
59	16.892	2359	2364	2374	rVB	207423	348710	5.56%	0.735%
60	17.069	2387	2394	2397	rBV	1375392	1895173	30.20%	3.992%

61	17.116	2397	2402	2410	rVV	3896660	6275498	100.00%	13.219%
62	17.186	2410	2414	2416	rVV2	180255	292058	4.65%	0.615%
63	17.227	2416	2421	2429	rVV2	449730	943300	15.03%	1.987%
64	17.316	2431	2436	2443	rVB2	64938	139270	2.22%	0.293%
65	17.492	2462	2466	2467	rBV4	9373	13362	0.21%	0.028%

66	17.522	2467	2471	2482	rVV	216402	467449	7.45%	0.985%
67	17.610	2482	2486	2489	rVB2	38456	52830	0.84%	0.111%
68	17.680	2495	2498	2501	rBV	18556	21899	0.35%	0.046%
69	17.827	2519	2523	2533	rVB4	24761	56240	0.90%	0.118%
70	17.980	2544	2549	2554	rBV	212856	364581	5.81%	0.768%

71	18.033	2554	2558	2569	rVV	230964	505080	8.05%	1.064%
72	18.133	2571	2575	2578	rVV	79232	125855	2.01%	0.265%
73	18.180	2578	2583	2588	rVV2	378176	787488	12.55%	1.659%
74	18.221	2588	2590	2598	rVB2	104001	167862	2.67%	0.354%
75	18.327	2605	2608	2613	rVB3	16284	24173	0.39%	0.051%

76	18.380	2613	2617	2624	rBV3	24229	54870	0.87%	0.116%
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 Title : SVOA CALIBRATION

77	18.510	2634	2639	2646	rVB	166097	229133	3.65%	0.483%
78	18.621	2656	2658	2664	rVB2	24064	30909	0.49%	0.065%
79	18.763	2677	2682	2688	rBV2	36847	64690	1.03%	0.136%
80	18.821	2688	2692	2695	rBV	33809	43877	0.70%	0.092%
81	18.863	2695	2699	2706	rBV3	24354	40228	0.64%	0.085%
82	18.939	2707	2712	2715	rBV	60962	93026	1.48%	0.196%
83	19.216	2752	2759	2768	rBV	2848120	4540345	72.35%	9.564%
84	19.468	2797	2802	2809	rBV2	106843	168823	2.69%	0.356%
85	19.592	2819	2823	2834	rVB	1991682	3550832	56.58%	7.480%
86	19.680	2835	2838	2844	rVB2	80886	115394	1.84%	0.243%
87	19.792	2853	2857	2862	rBV	119785	149373	2.38%	0.315%
88	19.880	2868	2872	2876	rBV	86764	129557	2.06%	0.273%
89	19.933	2878	2881	2890	rVB2	94968	221342	3.53%	0.466%
90	20.021	2892	2896	2901	rBV	83580	141912	2.26%	0.299%
91	20.121	2909	2913	2920	rVB	334041	487120	7.76%	1.026%
92	20.221	2925	2930	2934	rVV	388874	533101	8.49%	1.123%
93	20.274	2936	2939	2950	rVB3	128909	312720	4.98%	0.659%
94	21.092	3075	3078	3081	rBV	85219	110736	1.76%	0.233%
95	21.127	3081	3084	3088	rVB	118822	152706	2.43%	0.322%
96	21.510	3141	3149	3156	rBV2	1480634	4206673	67.03%	8.861%
97	21.715	3180	3184	3190	rVB	229977	361977	5.77%	0.762%
98	23.762	3525	3532	3538	rBV	477772	1353039	21.56%	2.850%
99	24.639	3675	3681	3692	rVB2	463075	1182214	18.84%	2.490%
100	24.792	3699	3707	3716	rBV	963036	2430546	38.73%	5.120%

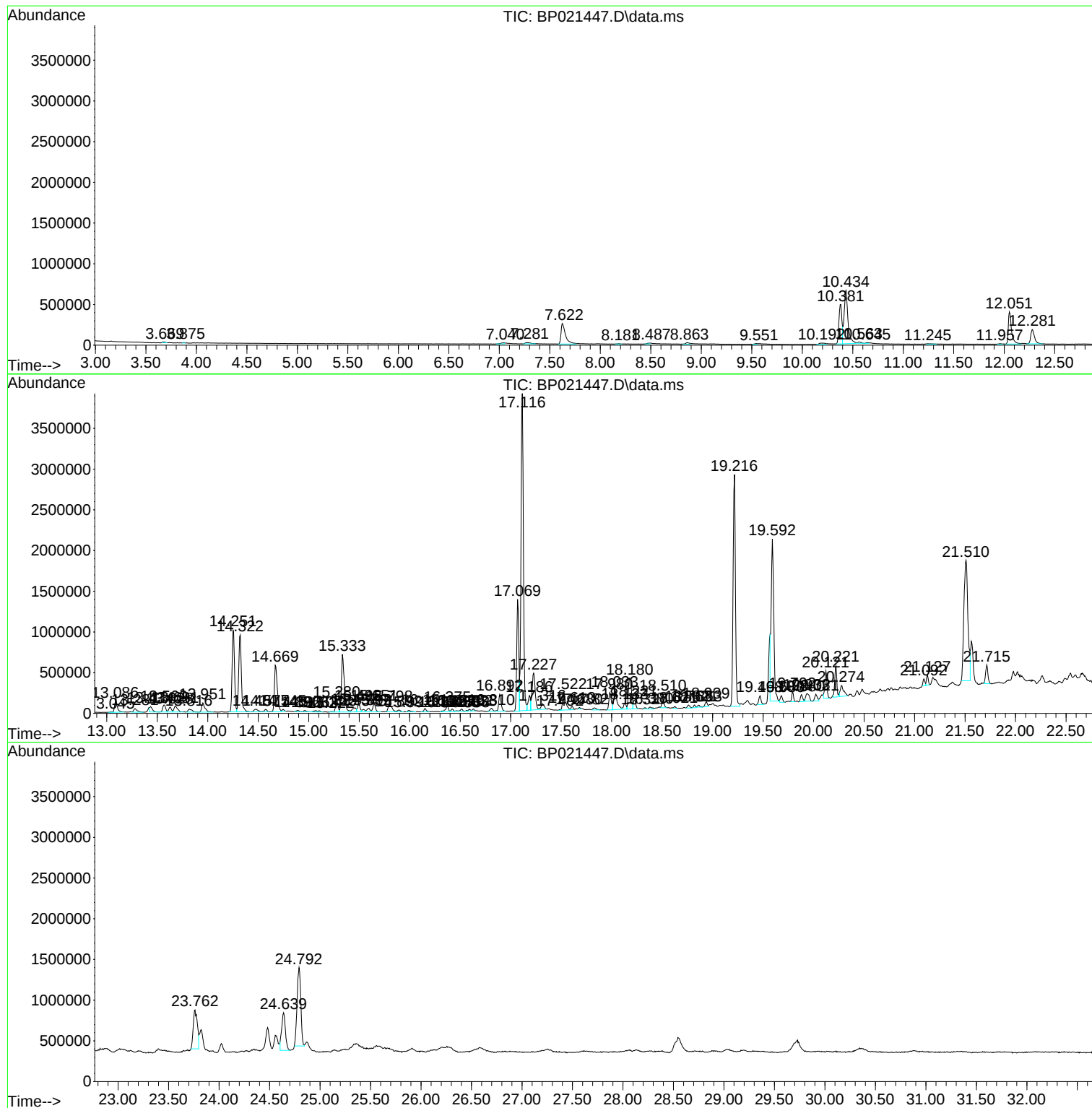
Sum of corrected areas: 47473178

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BNA_P
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Quant Title : SVOA CALIBRATION

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



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP080624\
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Misc :
ALS Vial : 39 Sample Multiplier: 1

Instrument :
BNA_P
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E05Y4DL

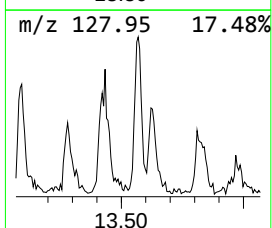
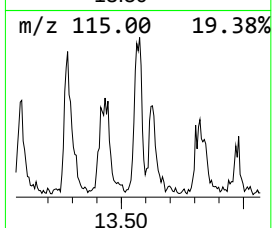
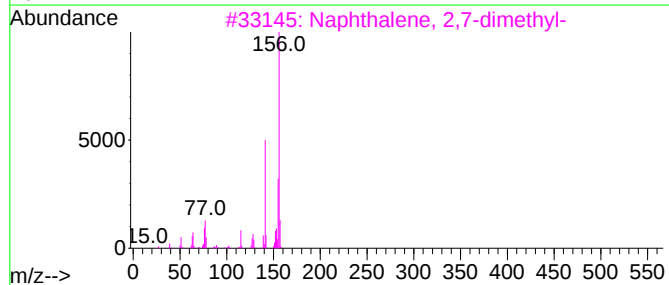
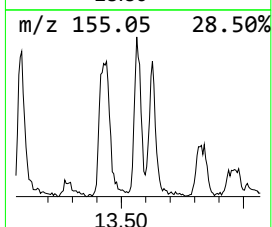
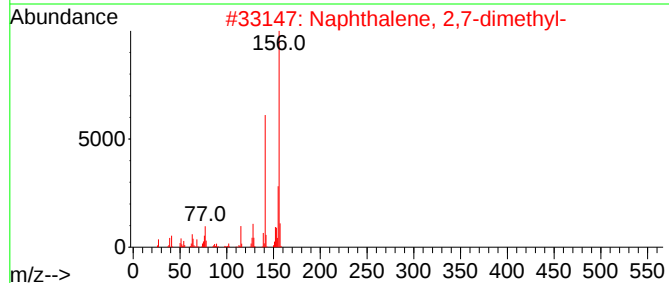
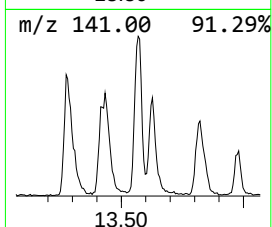
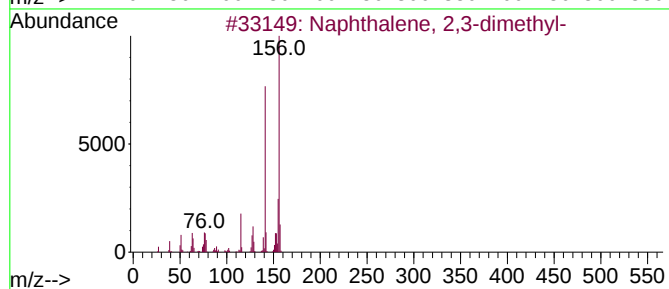
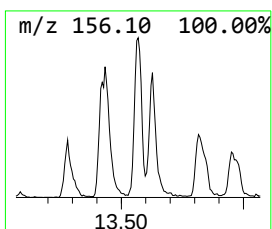
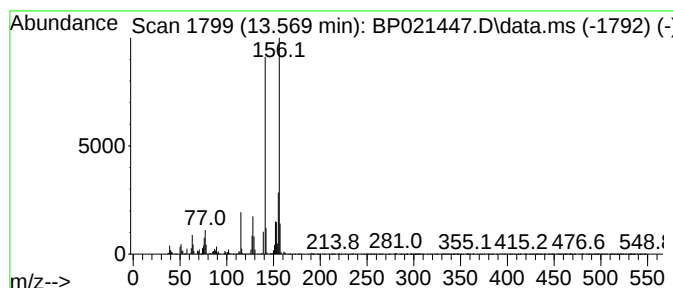
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Naphthalene, 2,3-dimethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.569	2.15 ng/ul	177592	Acenaphthene-d10	14.257

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	97
2			Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	97
3			Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	97
4			Naphthalene, 1,7-dimethyl-	156	C12H12	000575-37-1	97
5			Naphthalene, 1,3-dimethyl-	156	C12H12	000575-41-7	97



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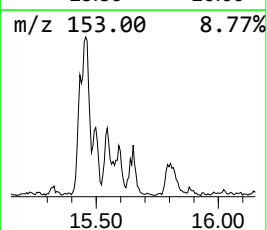
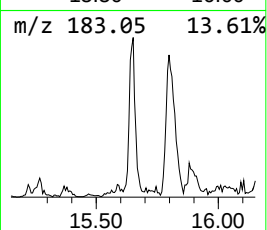
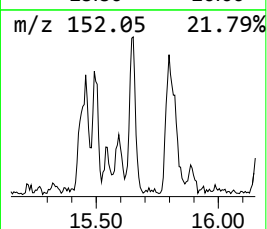
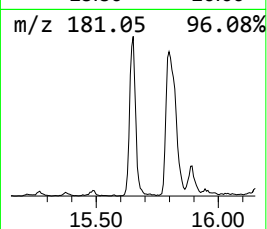
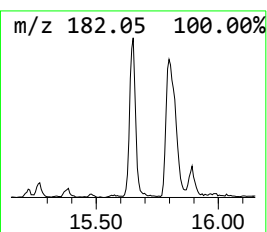
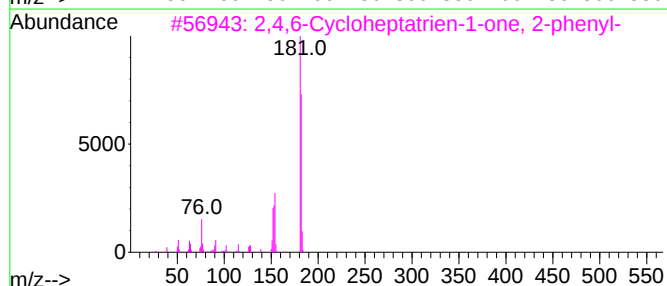
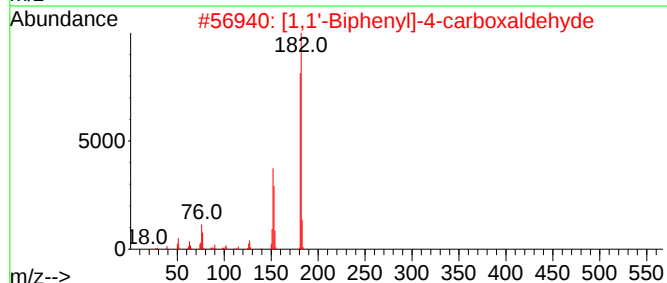
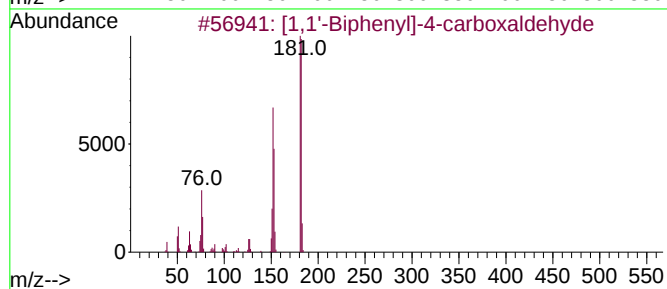
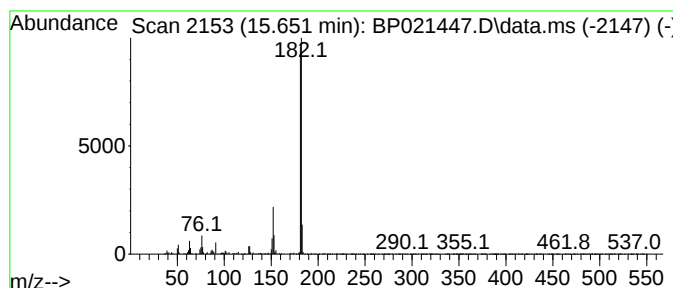
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Quant Title : SVOA CALIBRATION

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

Peak Number 2 [1,1'-Biphenyl]-4-carboxald... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.	
15.651	2.19 ng/ul	180848	Acenaphthene-d10	14.257	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	[1,1'-Biphenyl]-4-carboxaldehyde	182	C13H10O	003218-36-8	86
2	[1,1'-Biphenyl]-4-carboxaldehyde	182	C13H10O	003218-36-8	86
3	2,4,6-Cycloheptatrien-1-one, 2-p...	182	C13H10O	014562-09-5	83
4	9H-Fluoren-9-ol	182	C13H10O	001689-64-1	78
5	9H-Fluoren-9-ol	182	C13H10O	001689-64-1	78



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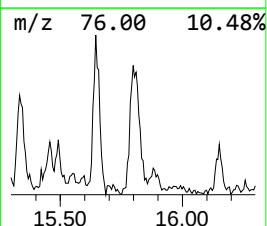
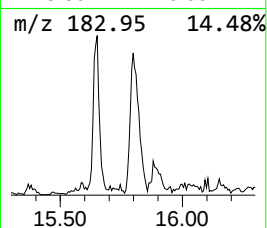
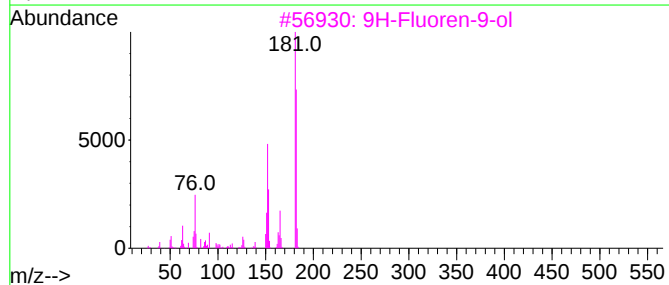
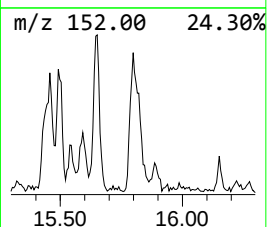
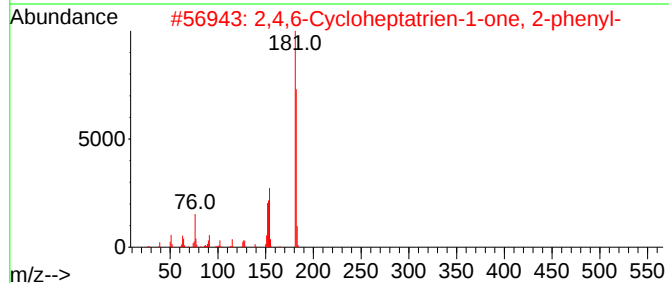
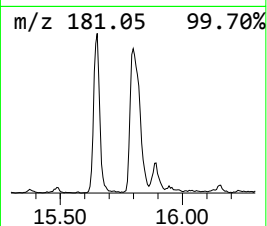
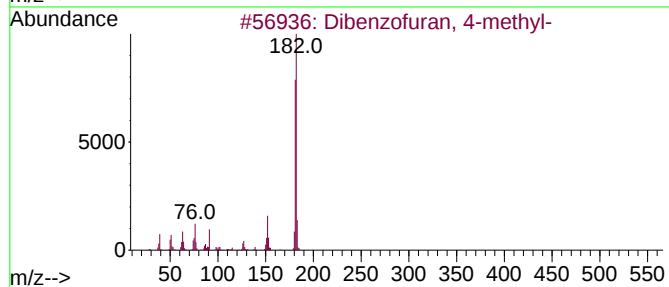
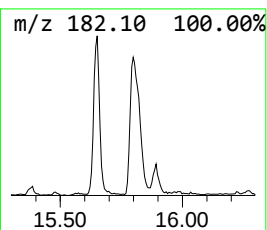
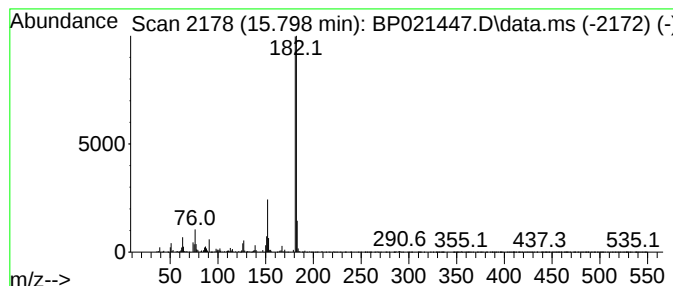
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 Dibenzofuran, 4-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.798	2.48 ng/ul	235475	Phenanthrene-d10	17.069

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Dibenzofuran, 4-methyl-	182	C13H10O	007320-53-8	91
2			2,4,6-Cycloheptatrien-1-one, 2-p...	182	C13H10O	014562-09-5	86
3			9H-Fluoren-9-ol	182	C13H10O	001689-64-1	78
4			9H-Fluoren-9-ol	182	C13H10O	001689-64-1	78
5			[1,1'-Biphenyl]-4-carboxaldehyde	182	C13H10O	003218-36-8	78



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Data File : BP021447.D
Acq On : 08 Aug 2024 12:38
Operator : CG/JU
Sample : P3437-14DL 10X
Misc :
ALS Vial : 39 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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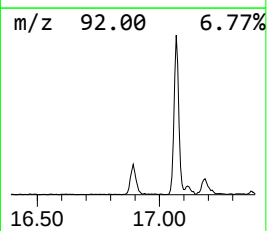
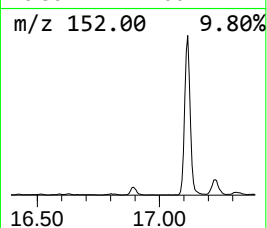
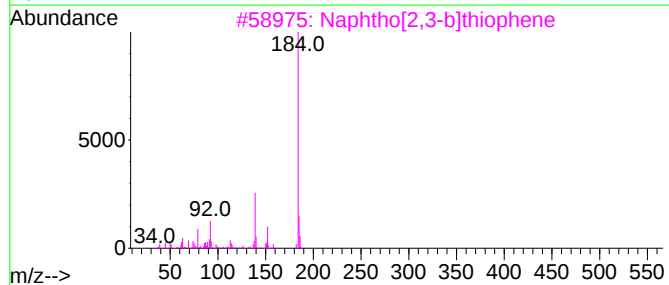
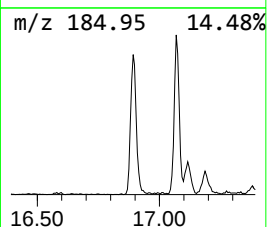
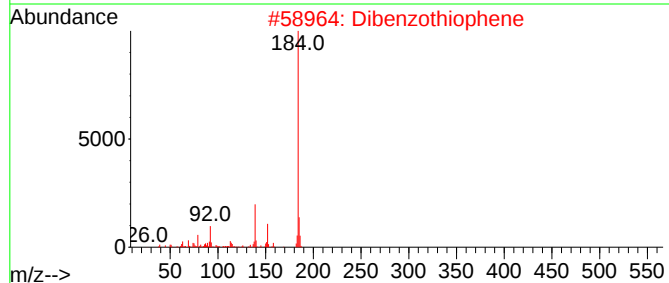
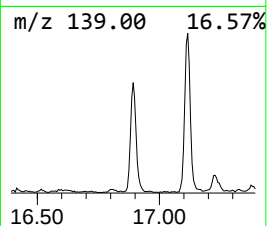
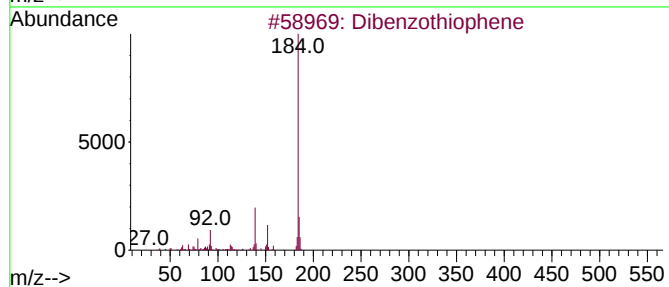
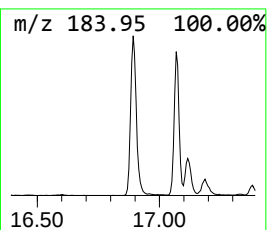
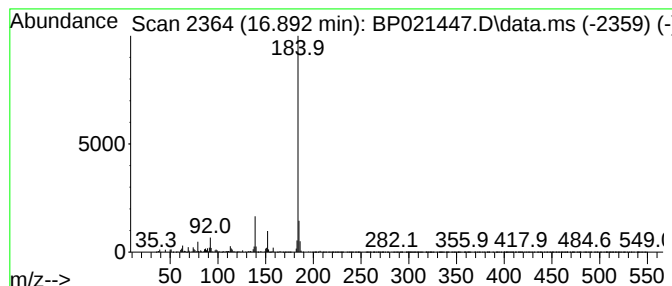
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 Dibenzothiophene Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.892	3.68 ng/ul	348710	Phenanthrene-d10	17.069

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Dibenzothiophene	184	C12H8S	000132-65-0	97
2			Dibenzothiophene	184	C12H8S	000132-65-0	97
3			Naphtho[2,3-b]thiophene	184	C12H8S	000268-77-9	97
4			Naphtho[2,1-b]thiophene	184	C12H8S	000233-02-3	96
5			Dibenzothiophene	184	C12H8S	000132-65-0	96



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ClientSampleId :
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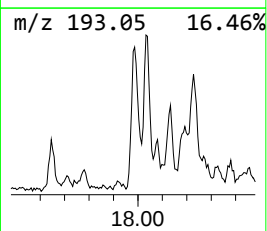
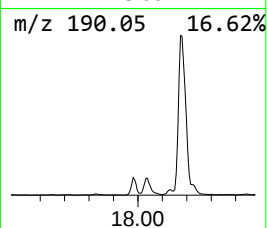
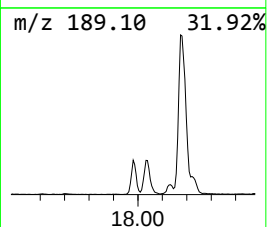
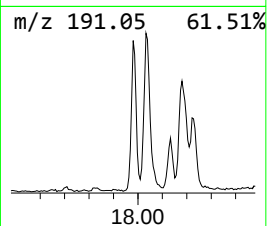
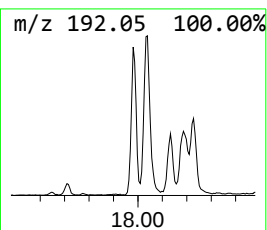
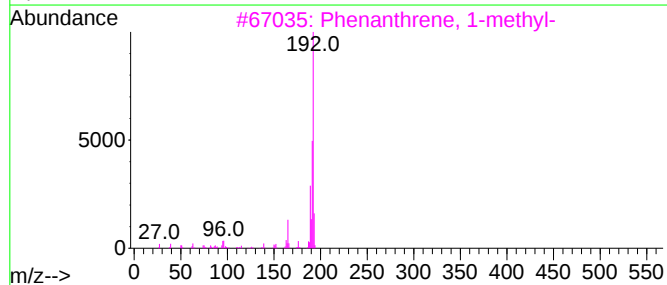
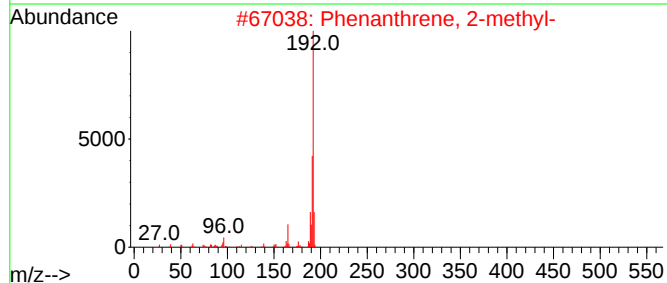
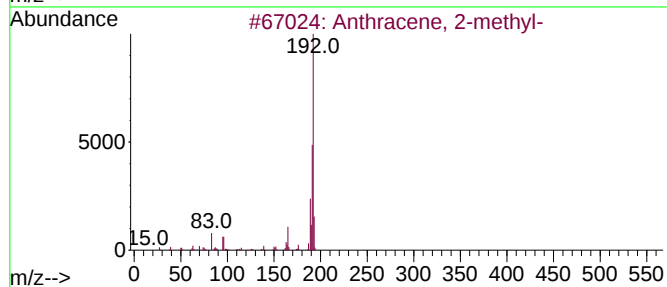
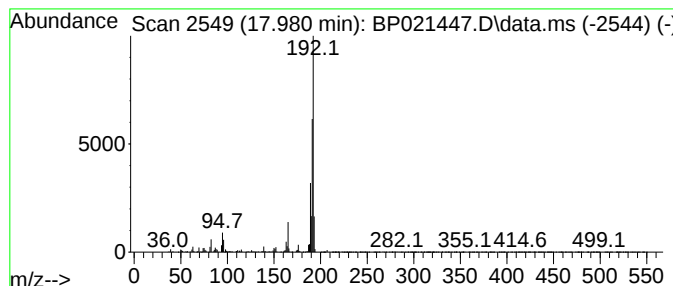
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 Anthracene, 2-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.980	3.85 ng/ul	364581	Phenanthrene-d10	17.069

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Anthracene, 2-methyl-	192	C15H12	000613-12-7	96
2			Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	96
3			Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	96
4			Phenanthrene, 4-methyl-	192	C15H12	000832-64-4	96
5			1H-Cyclopropa[1]phenanthrene,1a,...	192	C15H12	000949-41-7	95



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP080624\
Data File : BP021447.D
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Operator : CG/JU
Sample : P3437-14DL 10X
Misc :
ALS Vial : 39 Sample Multiplier: 1

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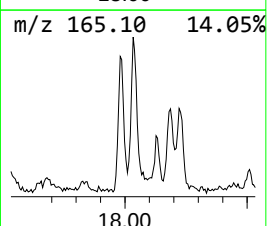
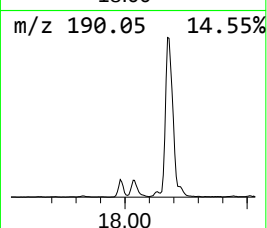
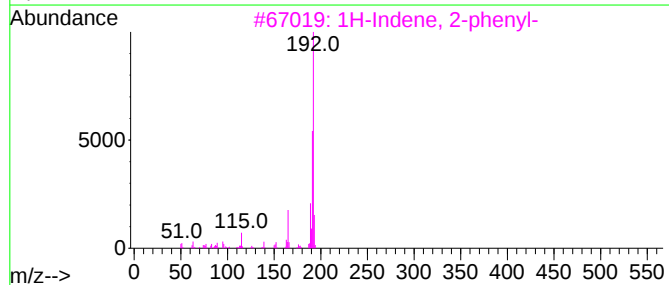
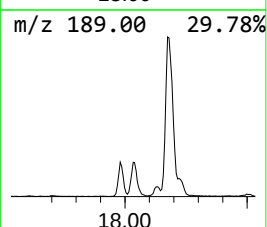
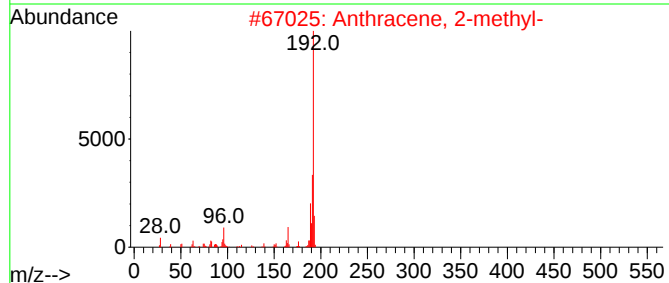
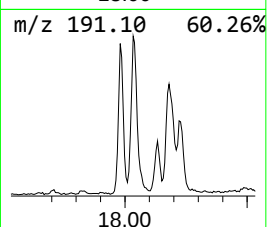
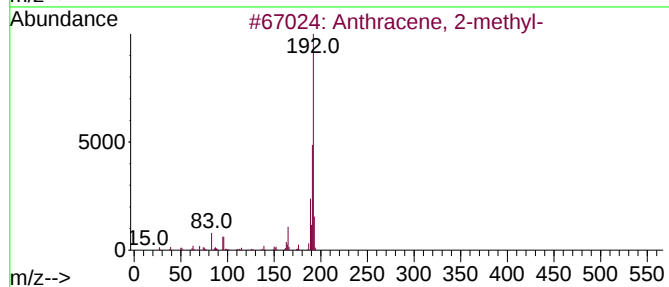
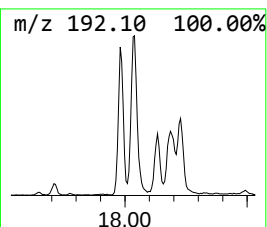
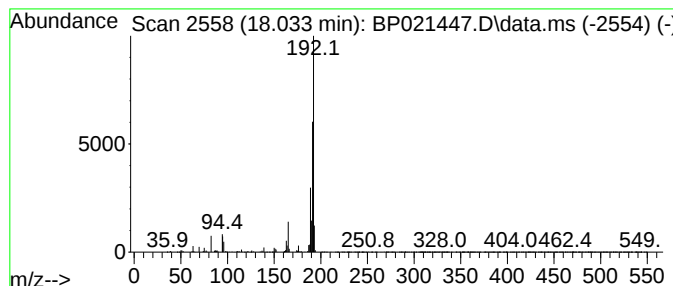
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 1H-Indene, 2-phenyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.033	5.33 ng/ul	505080	Phenanthrene-d10	17.069

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Anthracene, 2-methyl-	192	C15H12	000613-12-7	96
2		Anthracene, 2-methyl-	192	C15H12	000613-12-7	94
3		1H-Indene, 2-phenyl-	192	C15H12	004505-48-0	91
4		5H-Dibenzo[a,d]cycloheptene	192	C15H12	000256-81-5	91
5		Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP080624\
Data File : BP021447.D
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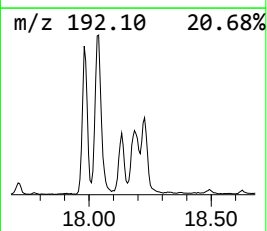
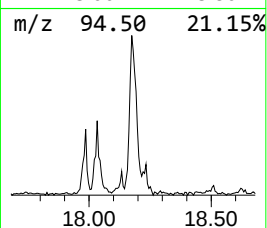
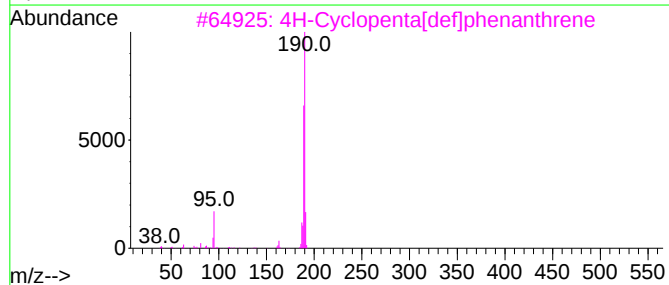
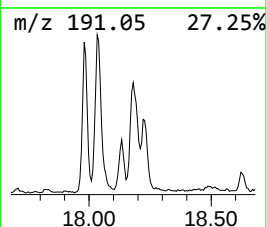
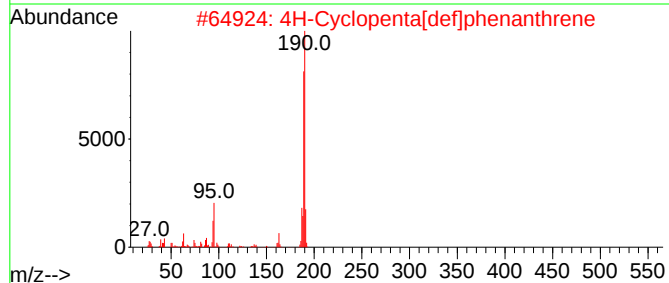
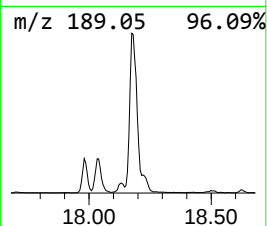
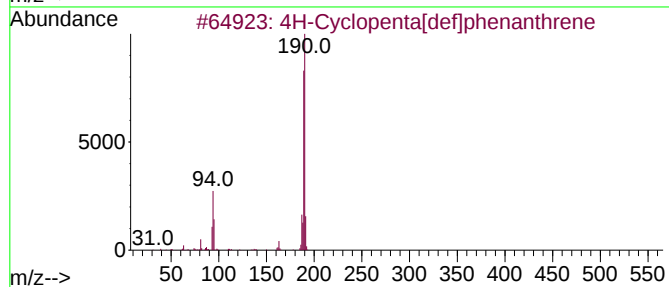
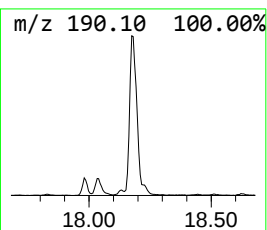
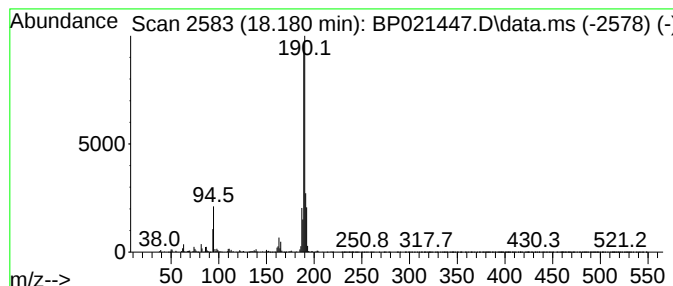
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 4H-Cyclopenta[def]phenanthrene Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.180	8.31 ng/ul	787488	Phenanthrene-d10	17.069

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	93
2			4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	76
3			4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	68
4			6H-Cyclobuta[jk]phenanthrene	190	C15H10	083469-43-6	64
5			2,3,5,6-Tetramethylterephthalald...	190	C12H14O2	007072-01-7	50



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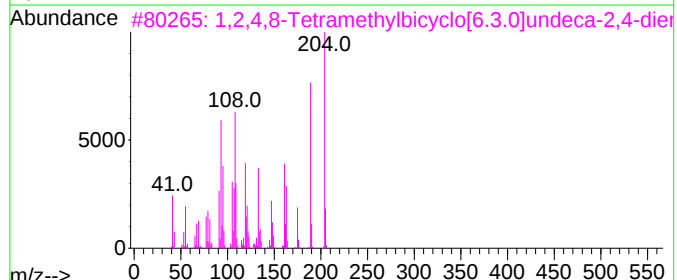
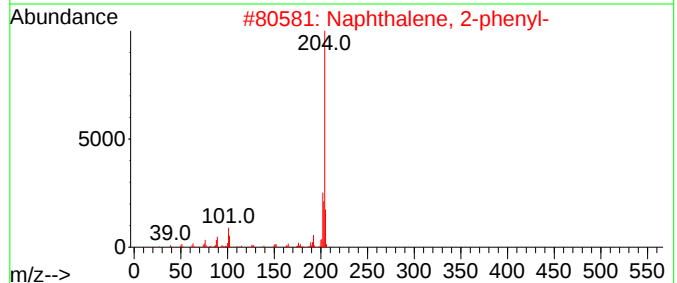
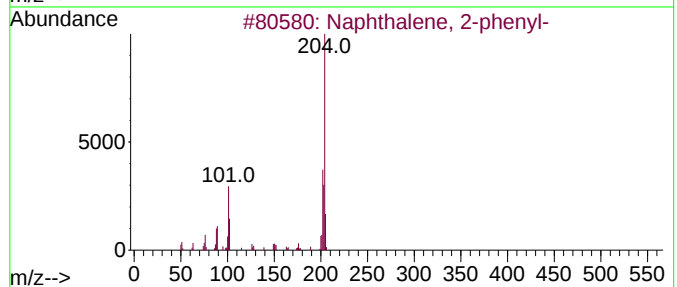
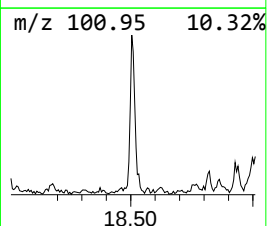
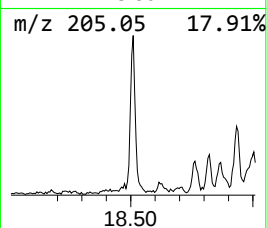
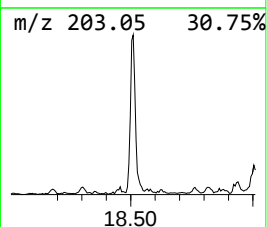
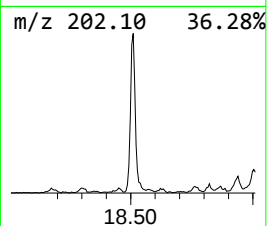
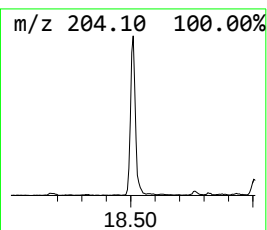
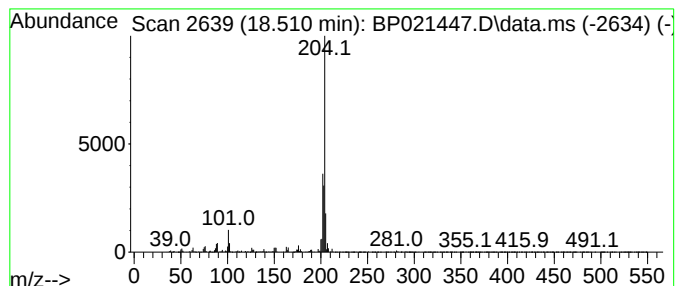
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 Naphthalene, 2-phenyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.510	2.42 ng/ul	229133	Phenanthrene-d10	17.069

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 2-phenyl-	204	C16H12	000612-94-2	93
2			Naphthalene, 2-phenyl-	204	C16H12	000612-94-2	93
3			1,2,4,8-Tetramethylbicyclo[6.3.0...	204	C15H24	137235-51-9	90
4			6-Phenylbenzocyclohepten-7-one	232	C17H12O	093327-56-1	86
5			5,16[1',2'] :8,13[1'',2'']-Dibenz...	408	C32H24	005672-97-9	86



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP080624\
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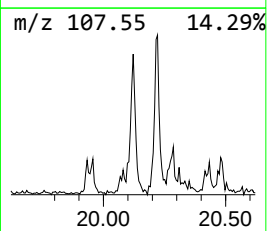
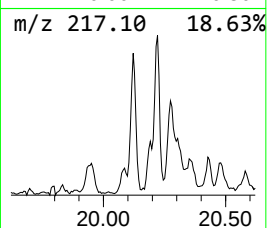
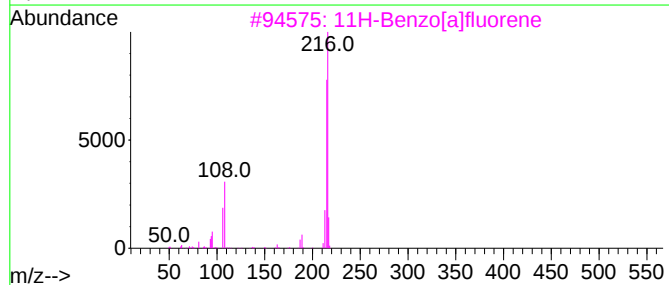
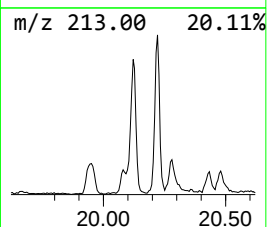
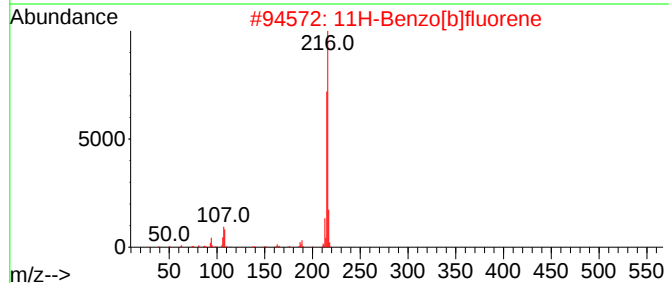
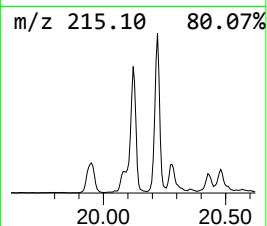
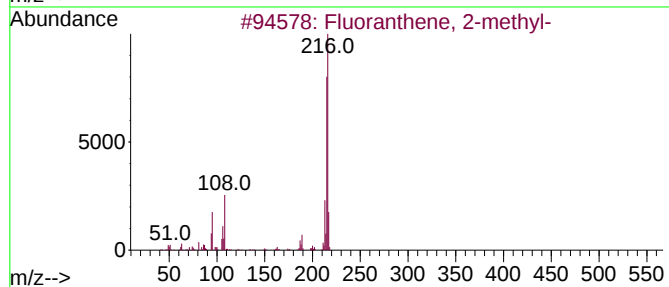
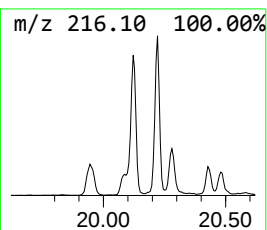
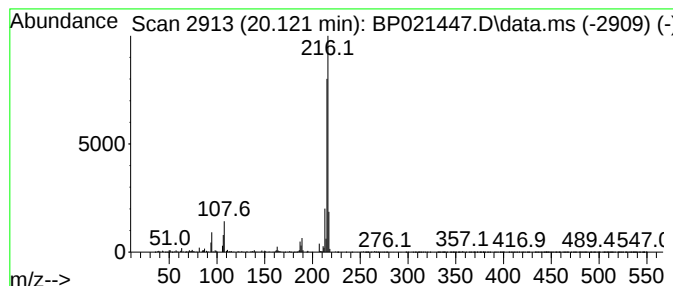
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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.121	2.32 ng/ul	487120	Chrysene-d12	21.515

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	95
2			11H-Benzo[b]fluorene	216	C17H12	000243-17-4	95
3			11H-Benzo[a]fluorene	216	C17H12	000238-84-6	91
4			11H-Benzo[b]fluorene	216	C17H12	000243-17-4	87
5			Pyrene, 1-methyl-	216	C17H12	002381-21-7	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP080624\
Data File : BP021447.D
Acq On : 08 Aug 2024 12:38
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ALS Vial : 39 Sample Multiplier: 1

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ClientSampleId :
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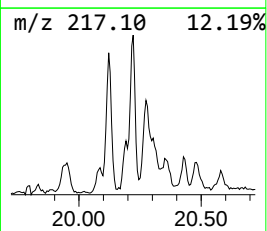
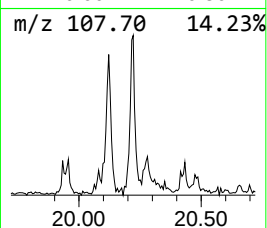
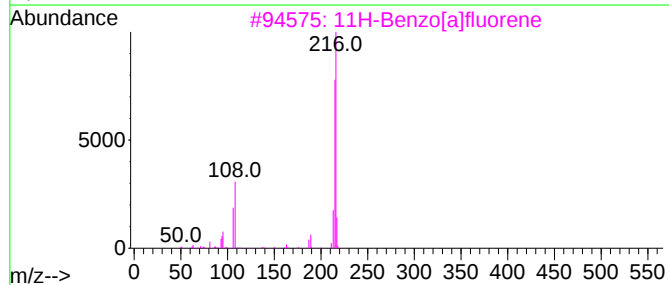
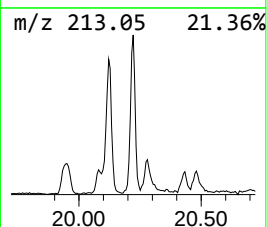
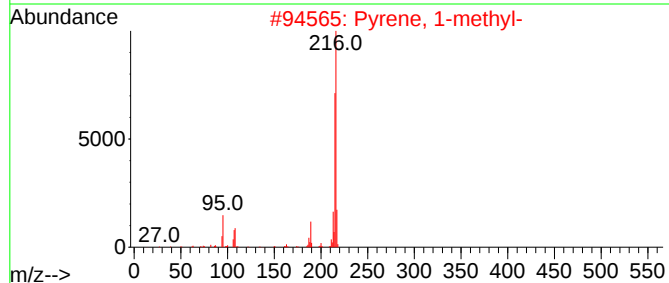
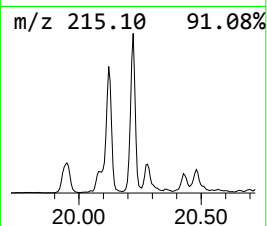
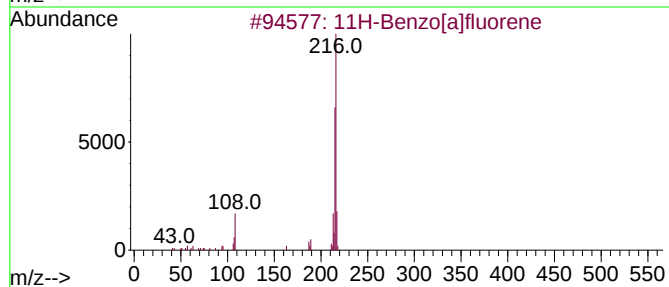
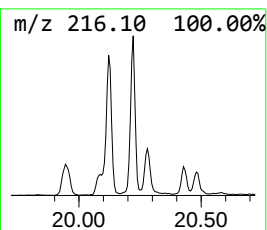
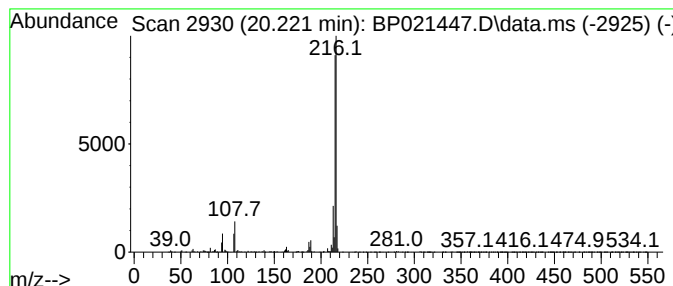
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 11H-Benzo[a]fluorene Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.221	2.53 ng/ul	533101	Chrysene-d12	21.515

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		11H-Benzo[a]fluorene	216	C17H12	000238-84-6	93
2		Pyrene, 1-methyl-	216	C17H12	002381-21-7	93
3		11H-Benzo[a]fluorene	216	C17H12	000238-84-6	90
4		11H-Benzo[b]fluorene	216	C17H12	000243-17-4	83
5		Pyrene, 1-methyl-	216	C17H12	002381-21-7	81



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP080624\
Data File : BP021447.D
Acq On : 08 Aug 2024 12:38
Operator : CG/JU
Sample : P3437-14DL 10X
Misc :
ALS Vial : 39 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
E05Y4DL

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.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
Naphthalene, 2,...	13.569	2.1	ng/ul	177592	3	14.257	1651660	20.0
[1,1'-Biphenyl]...	15.651	2.2	ng/ul	180848	3	14.257	1651660	20.0
Dibenzofuran, 4...	15.798	2.5	ng/ul	235475	4	17.069	1895170	20.0
Dibenzothiophene	16.892	3.7	ng/ul	348710	4	17.069	1895170	20.0
Anthracene, 2-m...	17.980	3.9	ng/ul	364581	4	17.069	1895170	20.0
1H-Indene, 2-ph...	18.033	5.3	ng/ul	505080	4	17.069	1895170	20.0
4H-Cyclopenta[d...	18.180	8.3	ng/ul	787488	4	17.069	1895170	20.0
Naphthalene, 2-...	18.510	2.4	ng/ul	229133	4	17.069	1895170	20.0
Fluoranthene, 2...	20.121	2.3	ng/ul	487120	5	21.515	4206670	20.0
11H-Benzo[a]flu...	20.221	2.5	ng/ul	533101	5	21.515	4206670	20.0