

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG060624\
 Data File : BG061522.D
 Acq On : 6 Jun 2024 23:13
 Operator : MA/JU
 Sample : P2634-03
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 BHBW9

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_G\Methods\SFAM-EPA-BG052124.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BG061522.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.111	12	21	42	rVB3	106712	405066	21.26%	2.019%
2	3.646	105	112	122	rBV2	7994	21319	1.12%	0.106%
3	3.781	126	135	148	rVV	41856	123742	6.49%	0.617%
4	6.566	605	609	615	rVB2	7563	12316	0.65%	0.061%
5	7.071	688	695	707	rVB	125428	225692	11.84%	1.125%
6	7.212	712	719	726	rBV	77407	130147	6.83%	0.649%
7	7.318	732	737	746	rBV3	13828	24499	1.29%	0.122%
8	7.424	748	755	761	rBV	115504	211040	11.07%	1.052%
9	7.882	826	833	839	rBV	125055	217025	11.39%	1.082%
10	8.599	948	955	968	rBB	146108	247062	12.97%	1.231%
11	9.034	1022	1029	1037	rBV	130734	227648	11.95%	1.135%
12	9.586	1118	1123	1128	rBV2	19099	27643	1.45%	0.138%
13	9.756	1143	1152	1159	rBV	125307	211287	11.09%	1.053%
14	10.309	1238	1246	1254	rBV	160942	291383	15.29%	1.452%
15	10.667	1300	1307	1313	rBV	155955	269446	14.14%	1.343%
16	10.808	1324	1331	1339	rBV	104832	196203	10.30%	0.978%
17	11.413	1424	1434	1445	rVB2	15351	33678	1.77%	0.168%
18	13.916	1850	1860	1866	rBV	288874	405963	21.30%	2.023%
19	14.204	1902	1909	1922	rVB	309004	505185	26.51%	2.518%
20	14.510	1951	1961	1968	rVB	242088	364424	19.12%	1.816%
21	14.756	1990	2003	2012	rBV2	72912	170687	8.96%	0.851%
22	15.503	2122	2130	2136	rBV	392467	593914	31.17%	2.960%
23	15.649	2150	2155	2163	rBV2	68646	109076	5.72%	0.544%
24	15.849	2184	2189	2195	rBV5	6051	11650	0.61%	0.058%
25	16.008	2205	2216	2222	rBV6	4265	11784	0.62%	0.059%
26	16.237	2249	2255	2260	rVB5	10678	21113	1.11%	0.105%
27	16.919	2366	2371	2377	rBV4	14471	20693	1.09%	0.103%
28	17.013	2381	2387	2391	rVB6	5489	12522	0.66%	0.062%
29	17.077	2395	2398	2405	rVB	11099	18863	0.99%	0.094%
30	17.259	2423	2429	2433	rBV	294538	453541	23.80%	2.261%
31	17.300	2433	2436	2441	rVV	190871	284693	14.94%	1.419%
32	17.359	2441	2446	2457	rVB	434688	664937	34.89%	3.314%
33	17.453	2457	2462	2467	rBV5	9256	18257	0.96%	0.091%
34	17.671	2490	2499	2503	rBV2	19453	40095	2.10%	0.200%
35	17.776	2514	2517	2520	rVB2	11306	11738	0.62%	0.059%
36	17.847	2524	2529	2536	rVB4	16220	30086	1.58%	0.150%

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Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG052124.MA.M
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37	18.064	2561	2566	2569	rBV6	7724	11955	0.63%	0.060%
38	18.135	2573	2578	2583	rVV	54644	116899	6.13%	0.583%
39	18.188	2583	2587	2591	rVV	78845	127666	6.70%	0.636%
40	18.270	2596	2601	2606	rVV2	12773	21695	1.14%	0.108%
41	18.329	2606	2611	2615	rVV2	70665	129688	6.81%	0.646%
42	18.370	2615	2618	2624	rVB2	39533	60388	3.17%	0.301%
43	18.452	2627	2632	2635	rBV7	10887	19204	1.01%	0.096%
44	18.534	2641	2646	2650	rVB6	9799	15338	0.80%	0.076%
45	18.634	2659	2663	2668	rVV	41907	65360	3.43%	0.326%
46	18.687	2668	2672	2679	rVB	53642	78608	4.13%	0.392%
47	18.857	2697	2701	2702	rBV4	9523	10989	0.58%	0.055%
48	18.887	2702	2706	2710	rVV	17252	25647	1.35%	0.128%
49	18.940	2710	2715	2717	rVV	20523	27355	1.44%	0.136%
50	18.975	2718	2721	2725	rVB4	17075	23176	1.22%	0.116%
51	19.057	2729	2735	2740	rBV2	55753	106441	5.59%	0.531%
52	19.116	2741	2745	2749	rVV4	22304	46421	2.44%	0.231%
53	19.145	2749	2750	2754	rVB	17572	16877	0.89%	0.084%
54	19.204	2755	2760	2766	rBV4	38071	77156	4.05%	0.385%
55	19.322	2771	2780	2786	rVV	594169	821105	43.09%	4.093%
56	19.380	2786	2790	2793	rVV2	19976	29256	1.54%	0.146%
57	19.421	2793	2797	2801	rVV3	21254	38646	2.03%	0.193%
58	19.469	2802	2805	2810	rVV	24182	40258	2.11%	0.201%
59	19.527	2812	2815	2817	rVV2	19715	29171	1.53%	0.145%
60	19.563	2819	2821	2824	rVV	28703	39248	2.06%	0.196%
61	19.592	2824	2826	2830	rVV4	18718	24010	1.26%	0.120%
62	19.657	2831	2837	2839	rVV2	630129	896176	47.03%	4.467%
63	19.686	2839	2842	2850	rVV	556958	777328	40.79%	3.874%
64	19.756	2850	2854	2860	rVB2	43394	71185	3.74%	0.355%
65	19.868	2868	2873	2879	rVB	24766	44305	2.32%	0.221%
66	19.921	2879	2882	2887	rVB4	28333	47615	2.50%	0.237%
67	20.003	2891	2896	2898	rBV4	55218	88159	4.63%	0.439%
68	20.144	2915	2920	2921	rBV4	40241	57754	3.03%	0.288%
69	20.179	2922	2926	2931	rVB	98978	159350	8.36%	0.794%
70	20.279	2939	2943	2946	rBV	54483	66024	3.46%	0.329%
71	20.338	2946	2953	2958	rVV2	70669	117740	6.18%	0.587%
72	20.473	2973	2976	2980	rBV	40424	52861	2.77%	0.263%
73	20.520	2981	2984	2987	rVV	48912	60772	3.19%	0.303%
74	20.573	2988	2993	2998	rVB	1575146	1905595	100.00%	9.498%
75	20.685	3010	3012	3015	rVB3	23546	23112	1.21%	0.115%
76	20.937	3051	3055	3060	rVB	70082	103536	5.43%	0.516%

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77	21.108	3080	3084	3088	rBV3	56229	87331	4.58%	0.435%
78	21.149	3088	3091	3093	rBV	38910	47534	2.49%	0.237%
79	21.261	3107	3110	3116	rVB	67642	99596	5.23%	0.496%
80	21.407	3131	3135	3140	rVB	147594	185174	9.72%	0.923%
81	21.513	3145	3153	3158	rVV3	473173	1180728	61.96%	5.885%
82	21.560	3158	3161	3172	rVB	309403	513663	26.96%	2.560%
83	21.707	3182	3186	3193	rBV2	48162	80681	4.23%	0.402%
84	21.778	3196	3198	3202	rVV2	21688	25613	1.34%	0.128%
85	21.913	3217	3221	3228	rVB3	36840	74119	3.89%	0.369%
86	22.224	3270	3274	3280	rVB2	105509	189914	9.97%	0.947%
87	22.283	3281	3284	3294	rVB3	54735	109630	5.75%	0.546%
88	22.383	3295	3301	3307	rBV5	62156	146826	7.70%	0.732%
89	22.488	3314	3319	3323	rBV4	81151	158667	8.33%	0.791%
90	22.870	3380	3384	3386	rBV5	36317	60910	3.20%	0.304%
91	22.923	3390	3393	3401	rVB4	73450	146185	7.67%	0.729%
92	23.282	3447	3454	3461	rBV6	44616	117192	6.15%	0.584%
93	23.434	3476	3480	3488	rVB5	100813	213790	11.22%	1.066%
94	23.634	3506	3514	3521	rBV	251734	763927	40.09%	3.808%
95	23.881	3552	3556	3564	rVB	55269	120325	6.31%	0.600%
96	24.328	3625	3632	3638	rBV	125894	330029	17.32%	1.645%
97	24.404	3638	3645	3651	rVV2	304412	816555	42.85%	4.070%
98	24.469	3651	3656	3666	rVB	217206	523400	27.47%	2.609%
99	24.610	3673	3680	3688	rBV	211494	541433	28.41%	2.699%
100	28.123	4267	4278	4297	rVB3	96519	431302	22.63%	2.150%

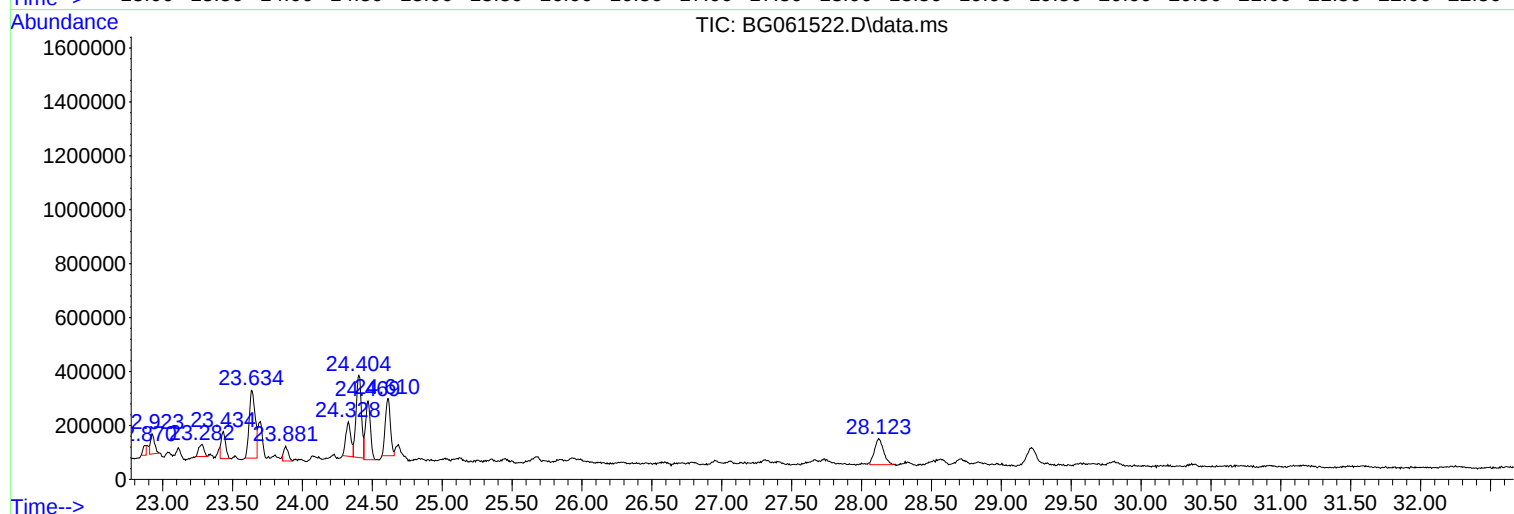
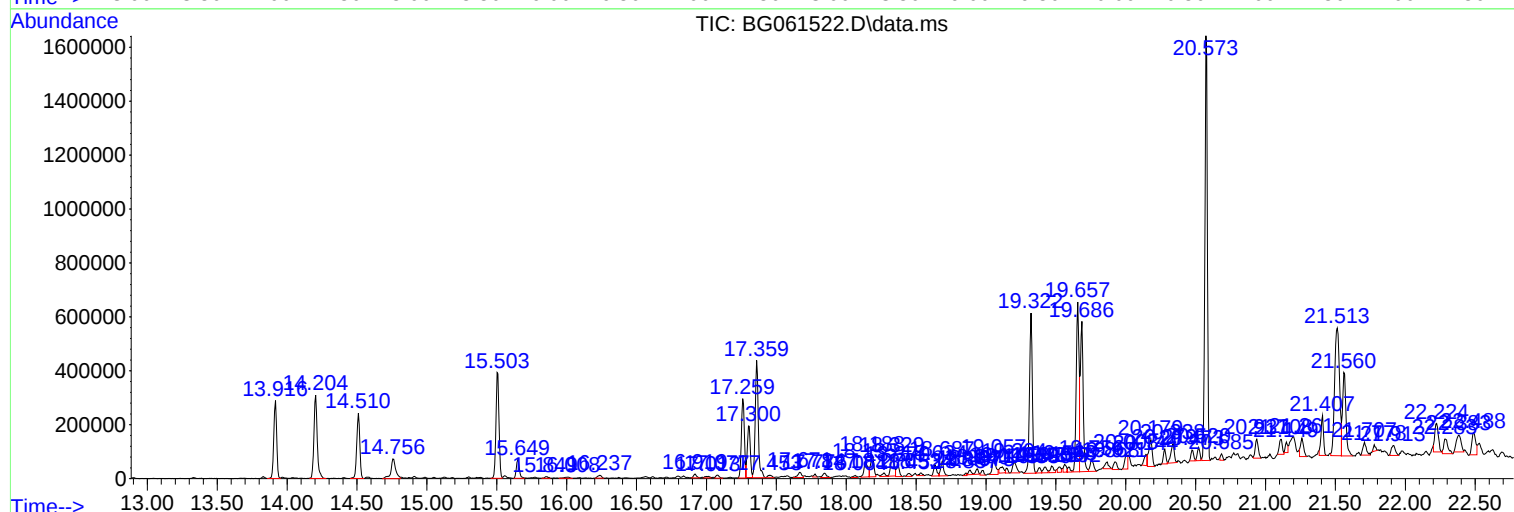
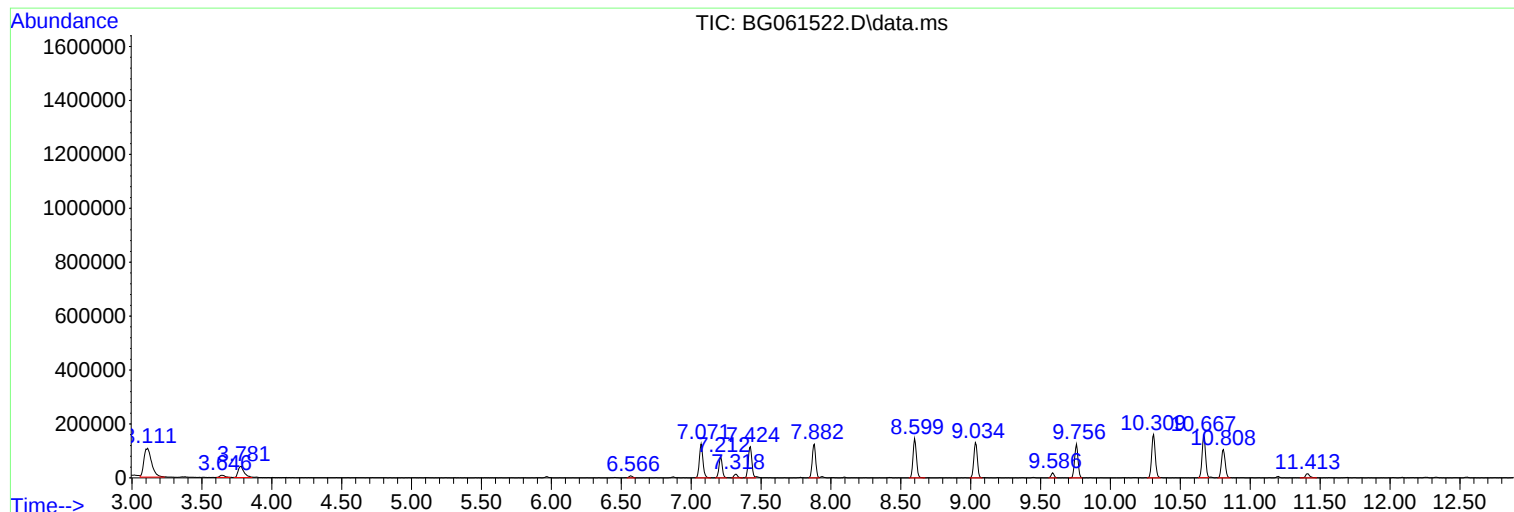
Sum of corrected areas: 20063010

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG052124.MA.M
Quant Title : SVOA CALIBRATION

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



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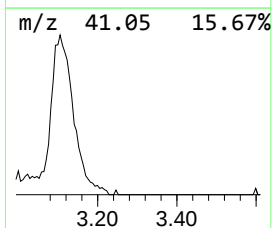
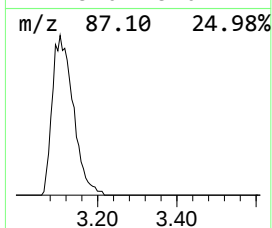
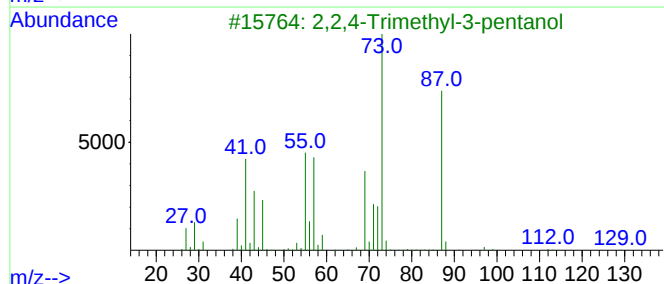
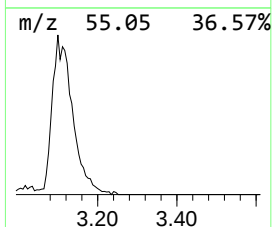
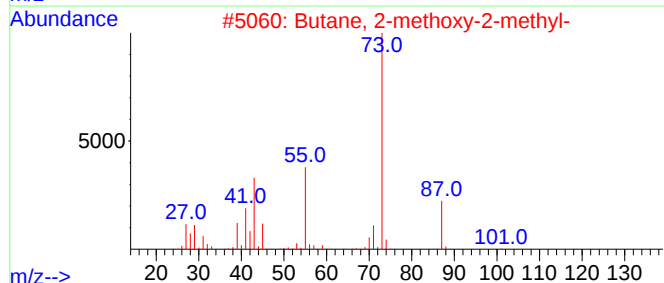
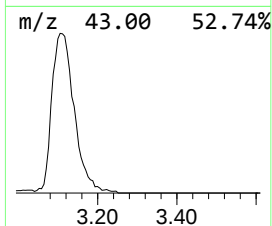
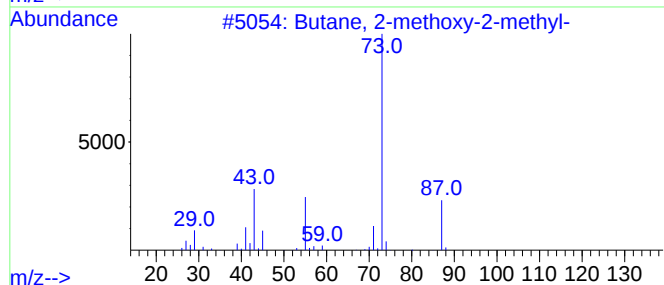
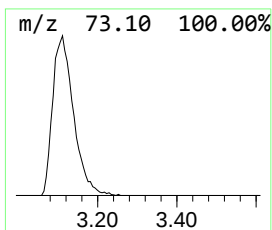
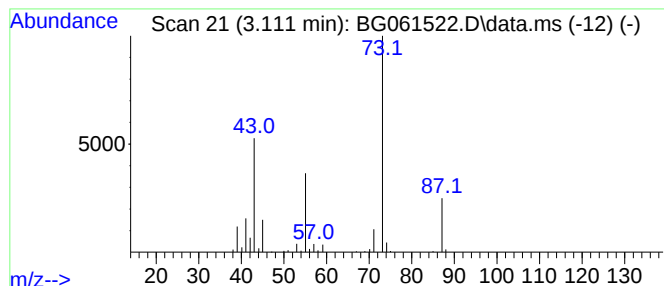
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG052124.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.111	37.33 ng/ul	405066	1,4-Dichlorobenzene-d4	7.882

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	83
2			Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	78
3			2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	39
4			2-Methyl-5-hexen-3-ol	114	C7H14O	032815-70-6	25
5			Silane, tetramethyl-	88	C4H12Si	000075-76-3	9



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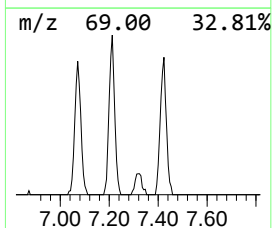
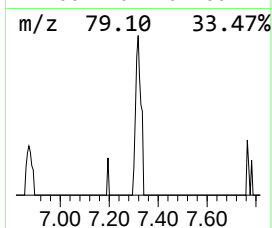
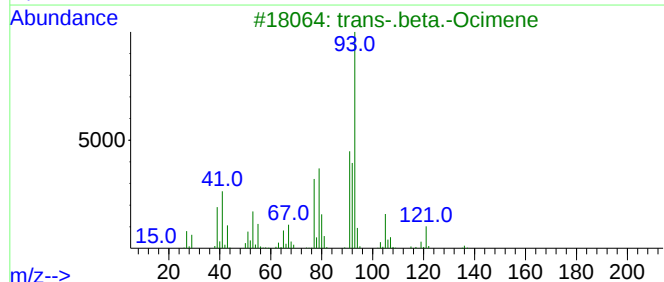
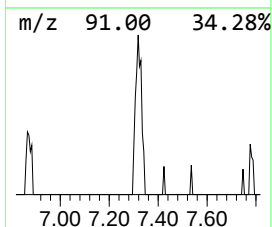
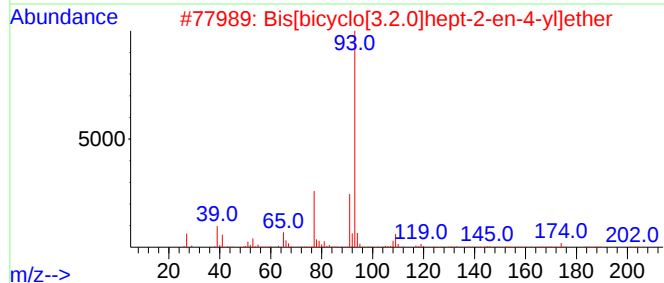
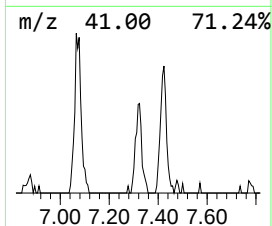
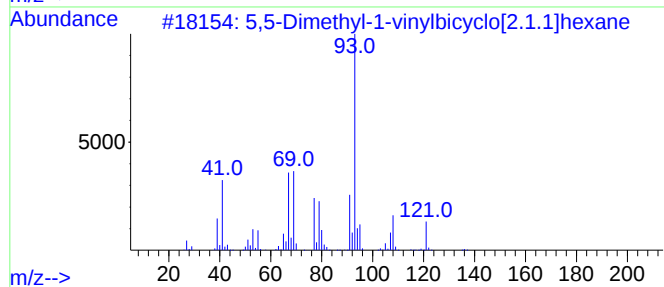
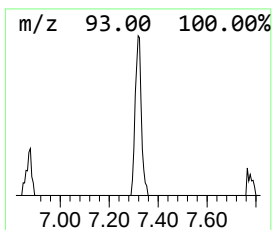
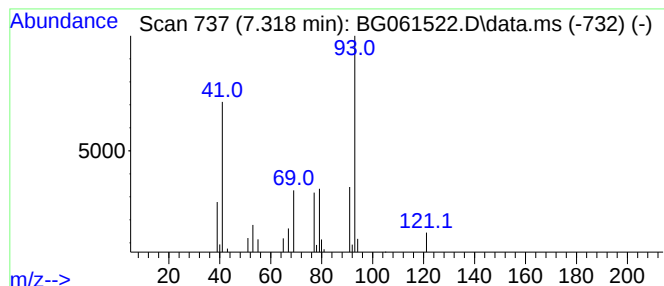
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG052124.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 5,5-Dimethyl-1-vinylbicyclo... Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.318	2.26 ng/ul	24499	1,4-Dichlorobenzene-d4	7.882

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			5,5-Dimethyl-1-vinylbicyclo[2.1.1]hexane	136	C10H16	016626-39-4	64
2			Bis[bicyclo[3.2.0]hept-2-en-4-yl]ether	202	C14H18O	1000153-89-5	59
3			trans-.beta.-Ocimene	136	C10H16	003779-61-1	56
4			3,5-Methanocyclopentapyrazole, 3,4-dihydro-2H-	164	C10H16N2	087143-58-6	45
5			trans-.beta.-Ocimene	136	C10H16	003779-61-1	45



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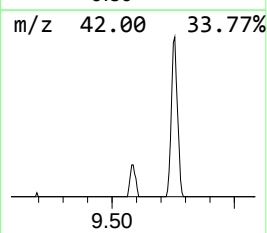
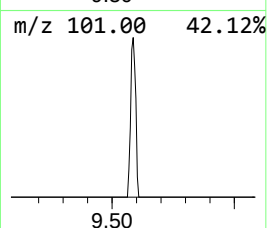
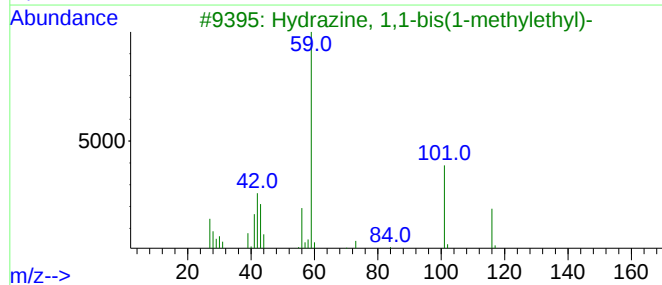
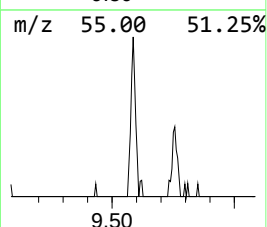
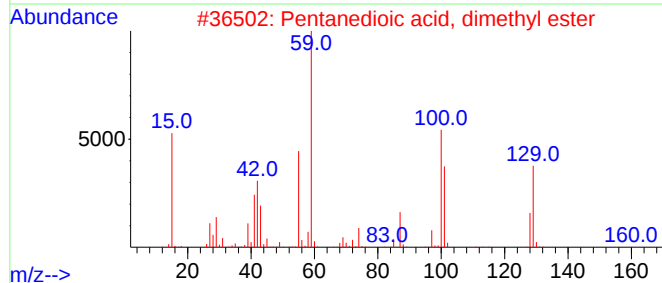
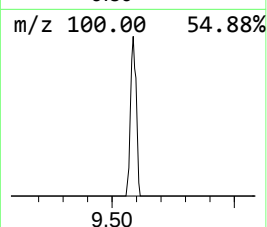
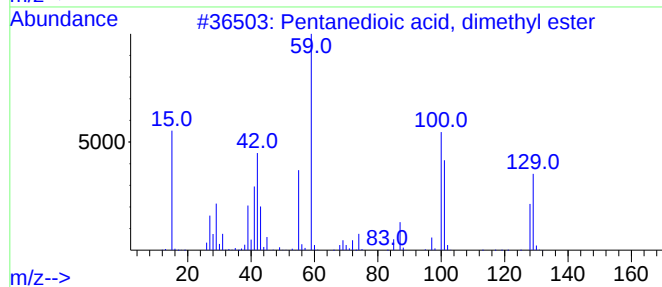
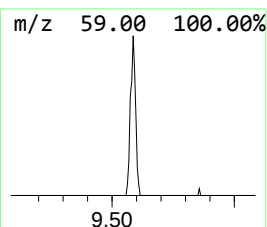
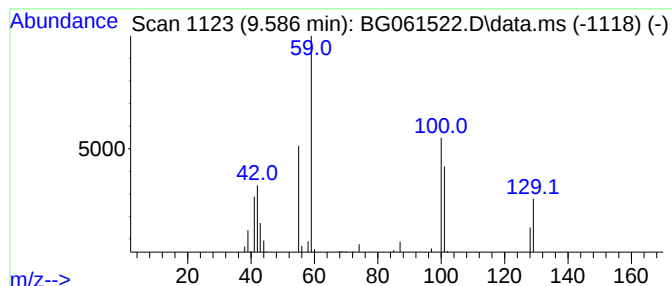
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Quant Title : SVOA CALIBRATION

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

Peak Number 3 Pentanedioic acid, dimethyl... Concentration Rank 21

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.586	2.05 ng/ul	27643	Naphthalene-d8	10.667

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentanedioic acid, dimethyl ester	160	C7H12O4	001119-40-0	90
2			Pentanedioic acid, dimethyl ester	160	C7H12O4	001119-40-0	43
3			Hydrazine, 1,1-bis(1-methylethyl)-	116	C6H16N2	000921-14-2	38
4			Propane, 2-methyl-2-(1-methylethyl)-	116	C7H16O	017348-59-3	37
5			Butanedioic acid, methyl-, dimethyl-	160	C7H12O4	001604-11-1	33



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG060624\
Data File : BG061522.D
Acq On : 6 Jun 2024 23:13
Operator : MA/JU
Sample : P2634-03
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BHBW9

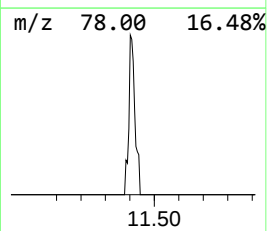
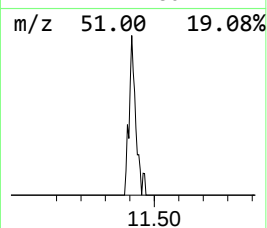
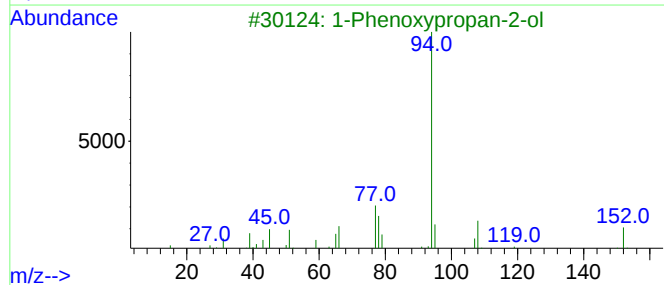
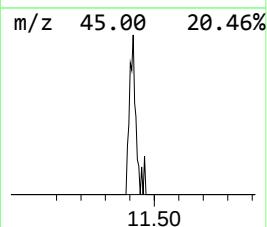
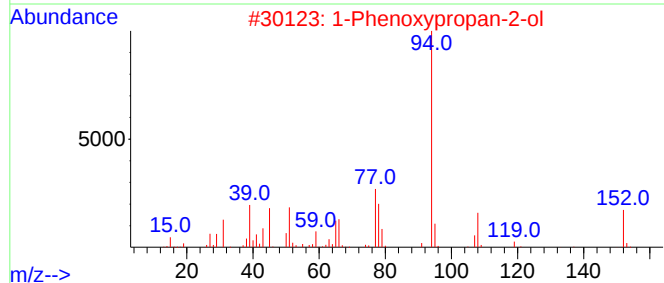
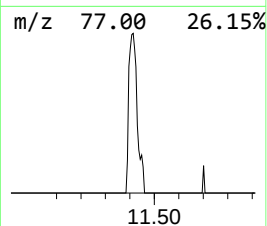
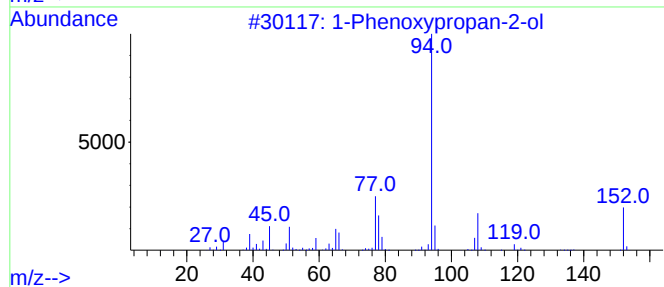
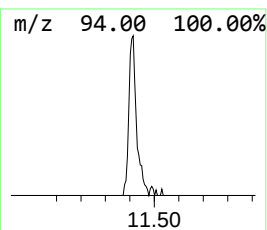
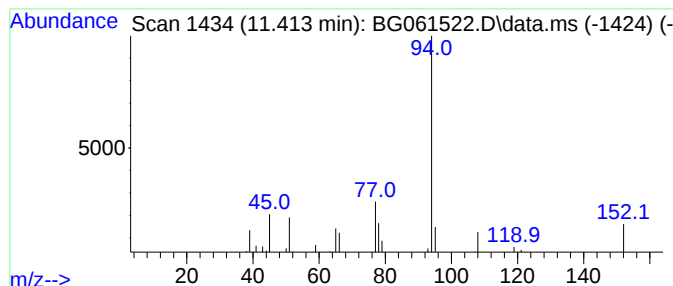
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 1-Phenoxypropan-2-ol Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.413	2.50 ng/ul	33678	Naphthalene-d8	10.667

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	94
2			1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	87
3			1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	86
4			1-Propanol, 3-phenoxy-	152	C9H12O2	006180-61-6	72
5			1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG060624\
Data File : BG061522.D
Acq On : 6 Jun 2024 23:13
Operator : MA/JU
Sample : P2634-03
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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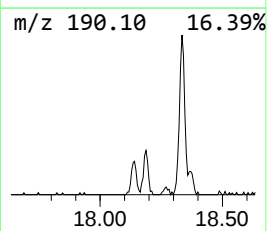
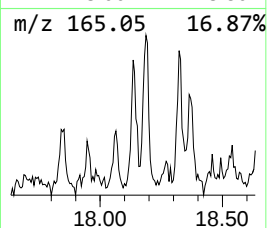
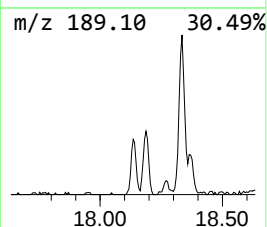
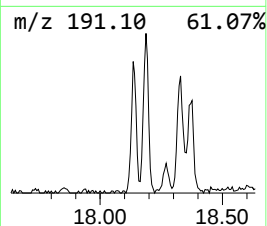
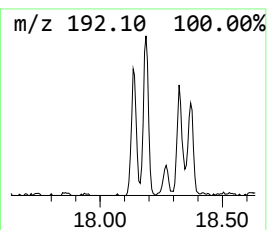
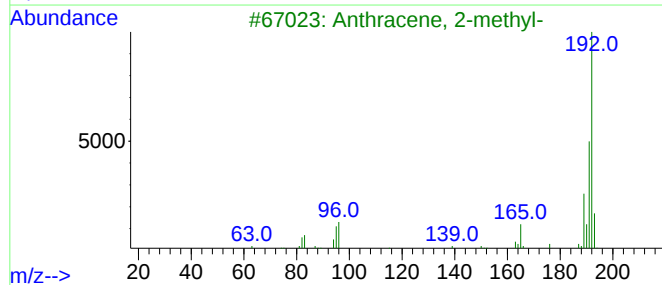
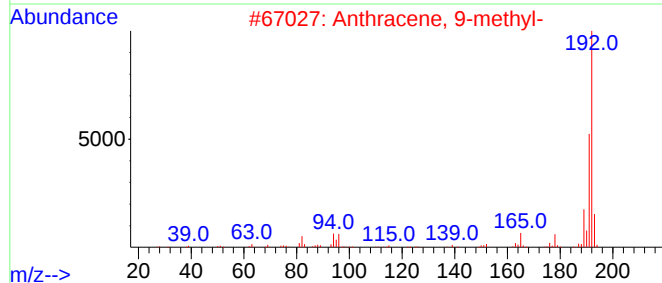
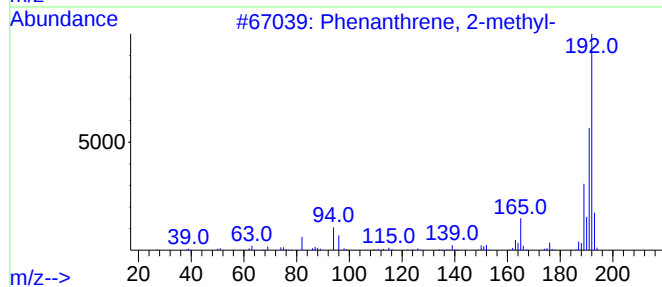
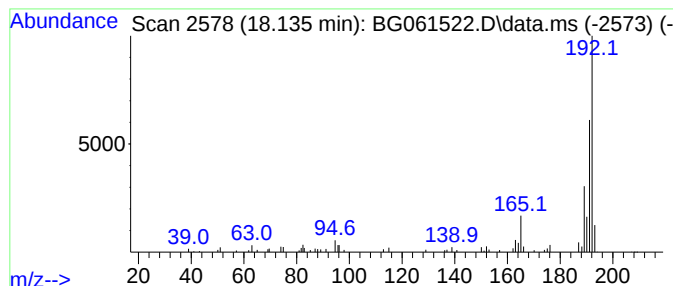
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TIC Integration Parameters: LSCINT.P

Peak Number 5 Phenanthrene, 2-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.135	5.15 ng/ul	116899	Phenanthrene-d10	17.259

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	94
2			Anthracene, 9-methyl-	192	C15H12	000779-02-2	90
3			Anthracene, 2-methyl-	192	C15H12	000613-12-7	90
4			Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	90
5			Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG060624\
Data File : BG061522.D
Acq On : 6 Jun 2024 23:13
Operator : MA/JU
Sample : P2634-03
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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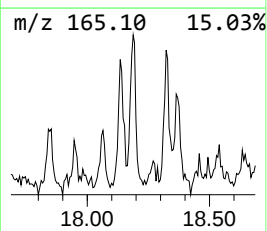
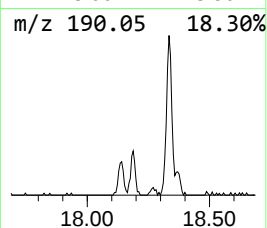
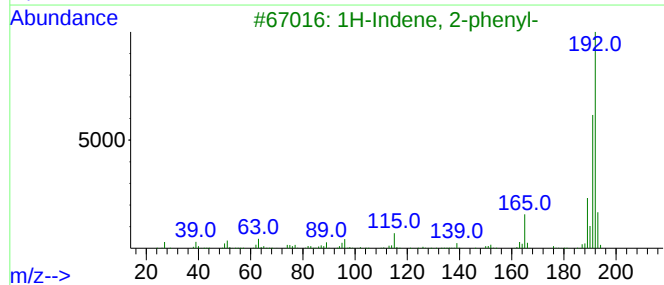
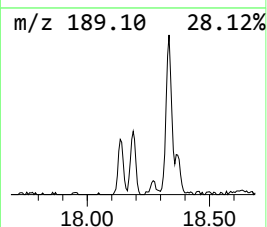
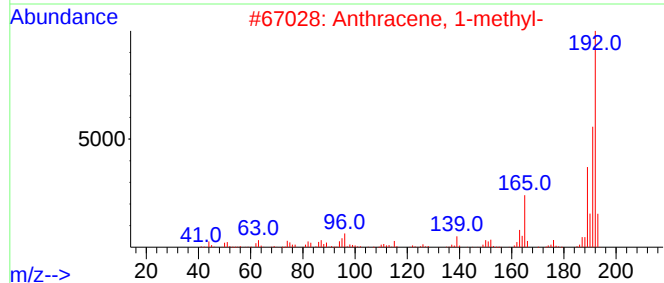
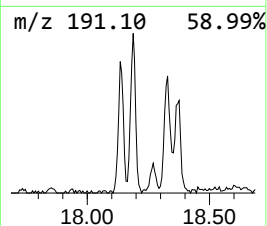
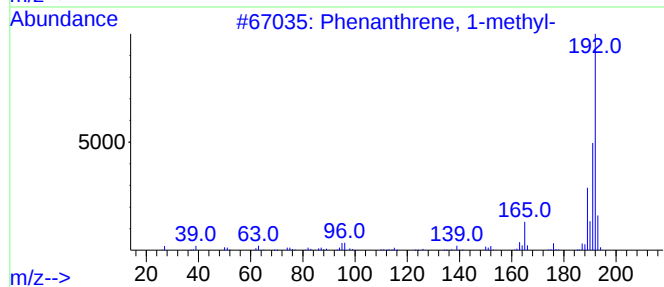
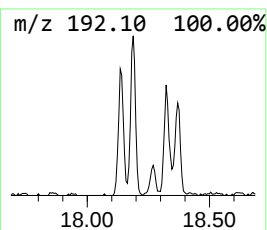
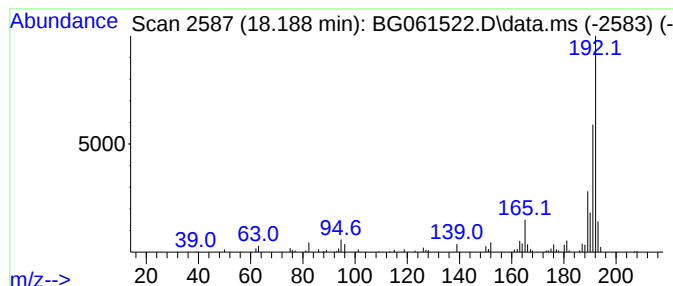
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 Phenanthrene, 1-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.188	5.63 ng/ul	127666	Phenanthrene-d10	17.259

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	96
2			Anthracene, 1-methyl-	192	C15H12	000610-48-0	93
3			1H-Indene, 2-phenyl-	192	C15H12	004505-48-0	93
4			Anthracene, 9-methyl-	192	C15H12	000779-02-2	93
5			Anthracene, 2-methyl-	192	C15H12	000613-12-7	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG060624\
Data File : BG061522.D
Acq On : 6 Jun 2024 23:13
Operator : MA/JU
Sample : P2634-03
Misc :
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Instrument :
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ClientSampleId :
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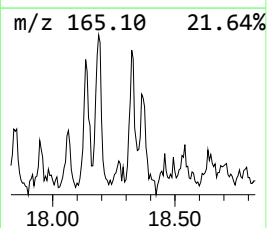
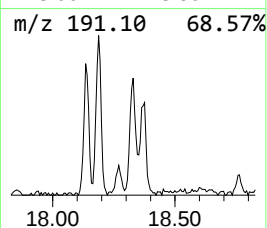
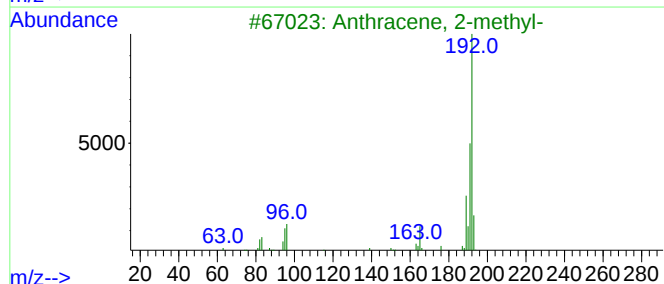
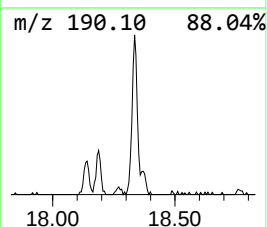
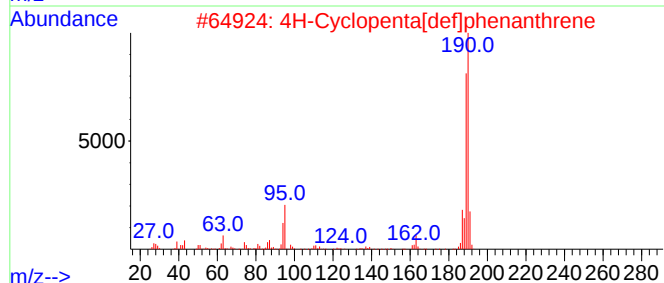
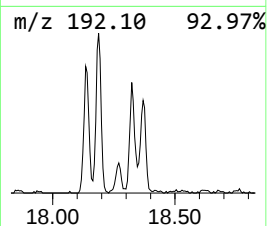
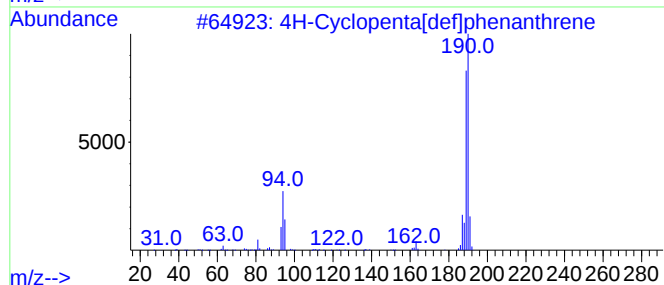
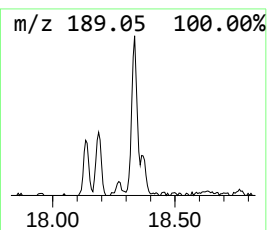
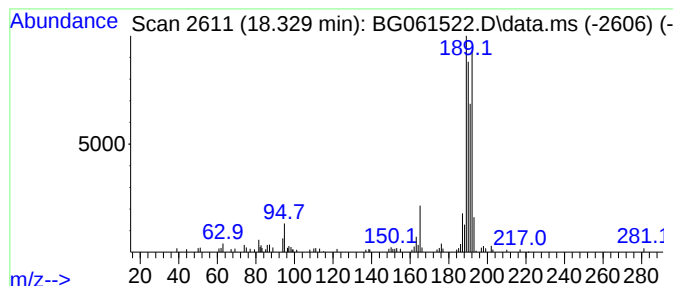
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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.329	5.72 ng/ul	129688	Phenanthrene-d10	17.259

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	60
2			4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	55
3			Anthracene, 2-methyl-	192	C15H12	000613-12-7	46
4			Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	45
5			Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	42



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG060624\
Data File : BG061522.D
Acq On : 6 Jun 2024 23:13
Operator : MA/JU
Sample : P2634-03
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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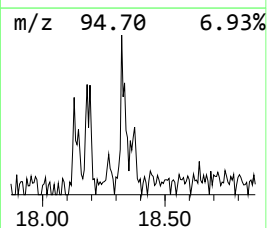
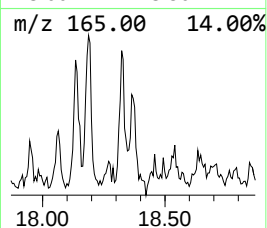
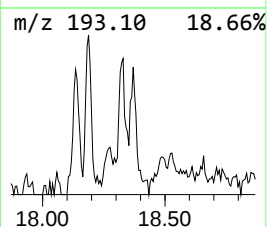
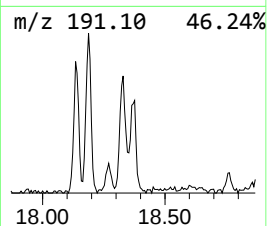
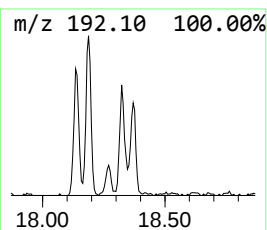
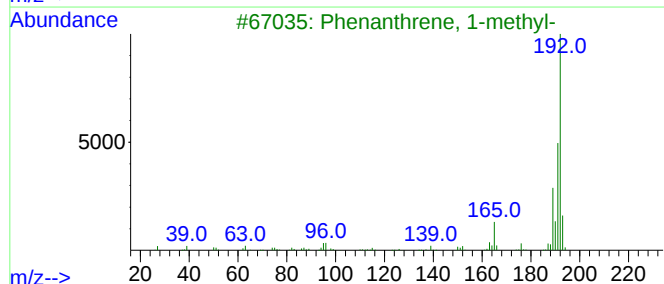
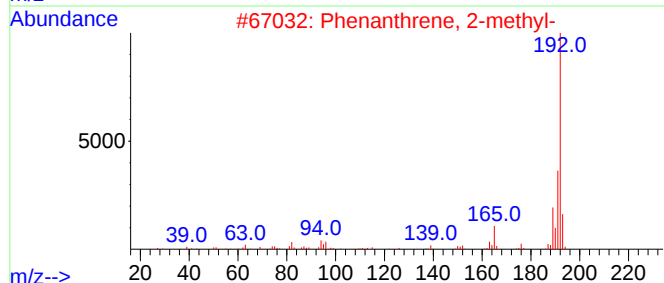
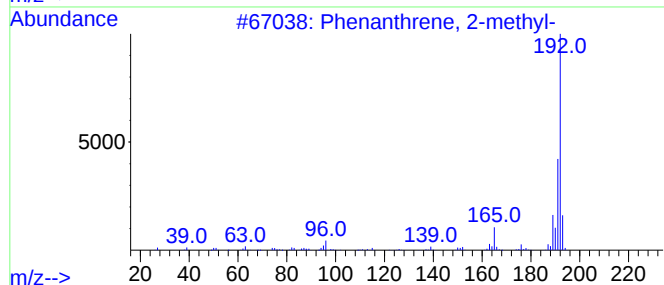
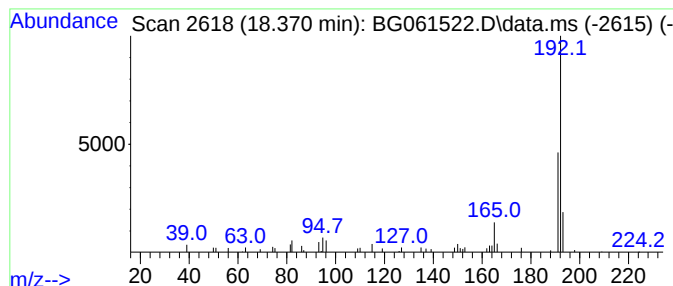
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 Phenanthrene, 3-methyl- Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.370	2.66 ng/ul	60388	Phenanthrene-d10	17.259

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	90
2			Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	90
3			Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	90
4			Phenanthrene, 3-methyl-	192	C15H12	000832-71-3	90
5			1H-Cyclopropa[1]phenanthrene,1a,...	192	C15H12	000949-41-7	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG060624\
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Acq On : 6 Jun 2024 23:13
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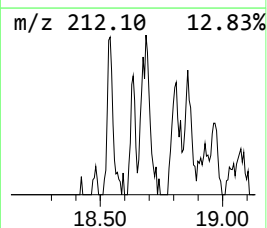
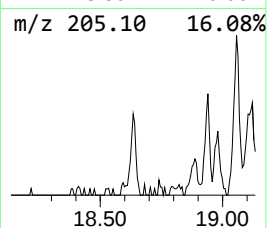
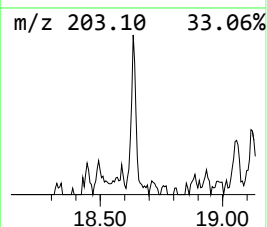
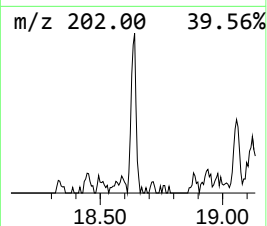
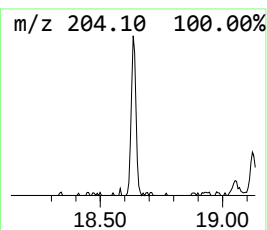
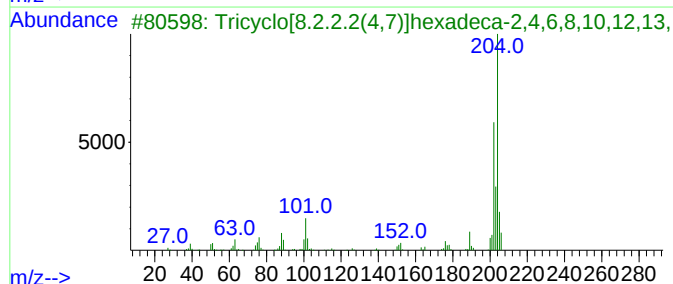
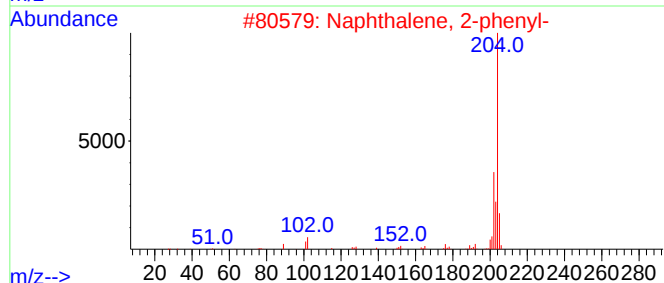
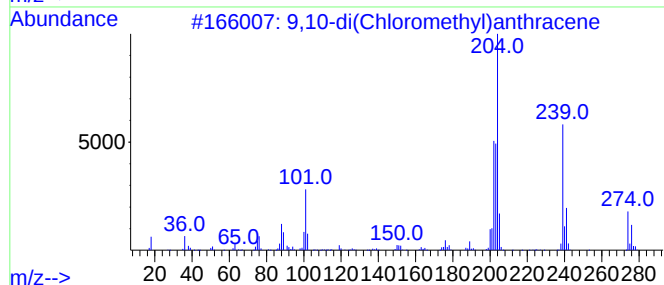
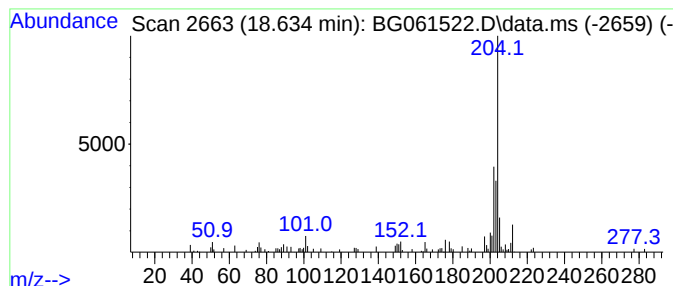
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 9,10-di(Chloromethyl)anthra... Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.634	2.88 ng/ul	65360	Phenanthrene-d10	17.259

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			9,10-di(Chloromethyl)anthracene	274	C16H12Cl2	010387-13-0	72
2			Naphthalene, 2-phenyl-	204	C16H12	000612-94-2	64
3			Tricyclo[8.2.2.2(4,7)]hexadeca-2...	204	C16H12	006572-60-7	58
4			9,10-Bis(bromomethyl)anthracene	362	C16H12Br2	034373-96-1	50
5			[1,1'-Biphenyl]-4,4'-dicarbonitrile	204	C14H8N2	001591-30-6	49



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG060624\
Data File : BG061522.D
Acq On : 6 Jun 2024 23:13
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Misc :
ALS Vial : 10 Sample Multiplier: 1

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ClientSampleId :
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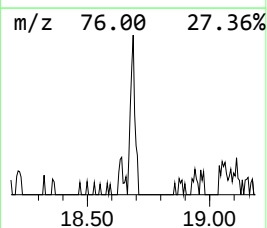
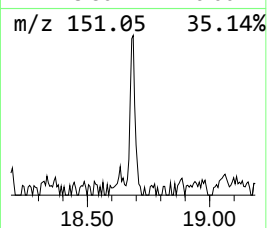
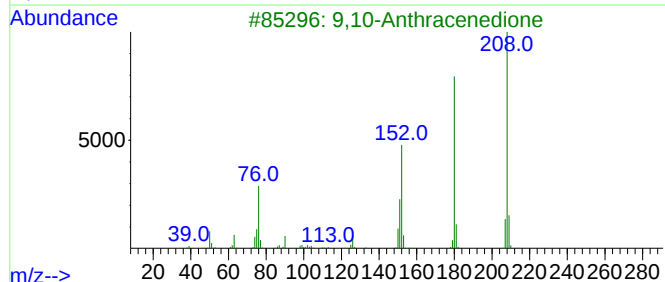
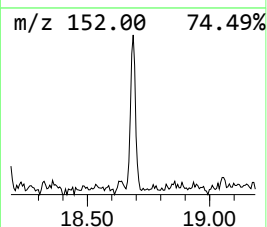
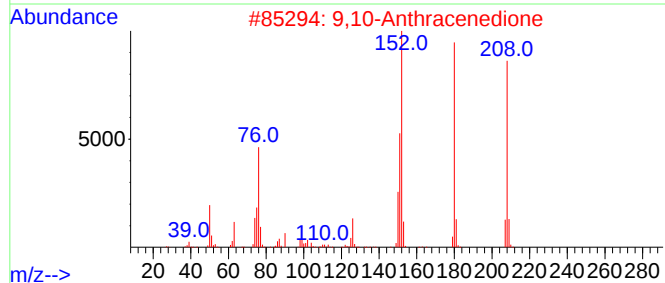
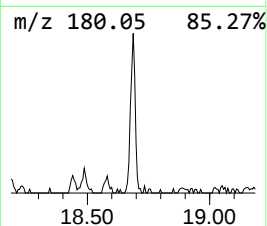
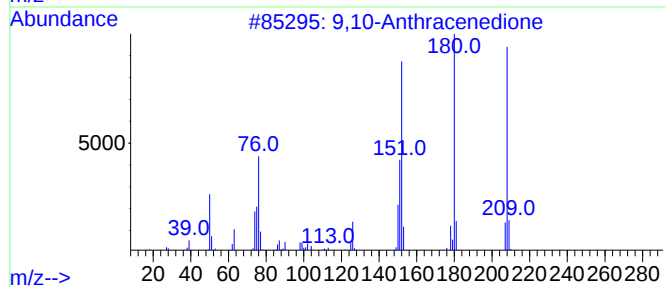
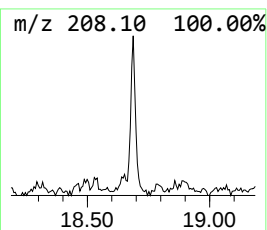
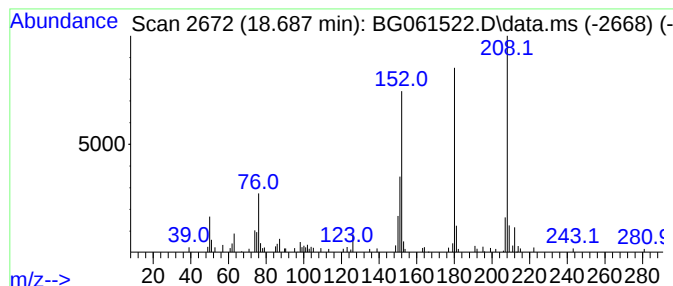
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 9,10-Anthracenedione Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.687	3.47 ng/ul	78608	Phenanthrene-d10	17.259

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	9,10-Anthracenedione			208	C14H8O2	000084-65-1	94
2	9,10-Anthracenedione			208	C14H8O2	000084-65-1	93
3	9,10-Anthracenedione			208	C14H8O2	000084-65-1	90
4	9,10-Anthracenedione			208	C14H8O2	000084-65-1	90
5	Benzo[c]cinnoline			180	C12H8N2	000230-17-1	78



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG060624\
Data File : BG061522.D
Acq On : 6 Jun 2024 23:13
Operator : MA/JU
Sample : P2634-03
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BHBW9

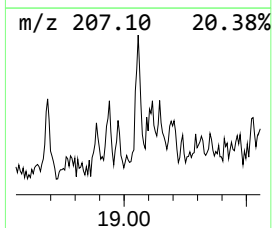
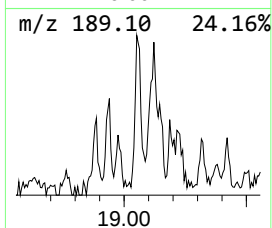
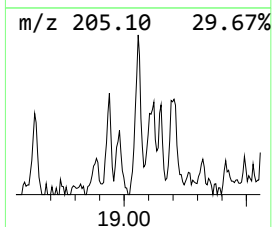
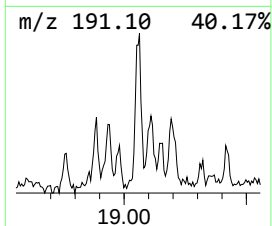
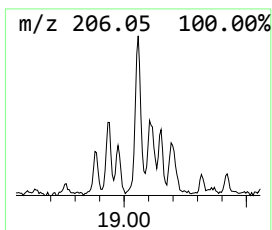
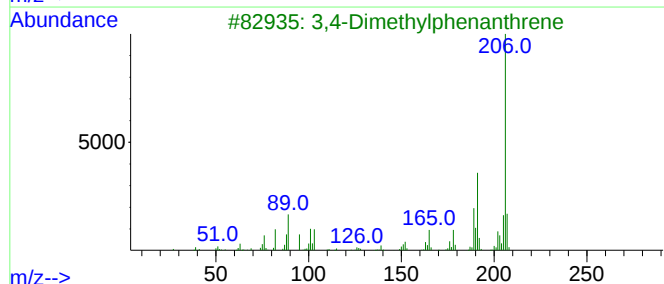
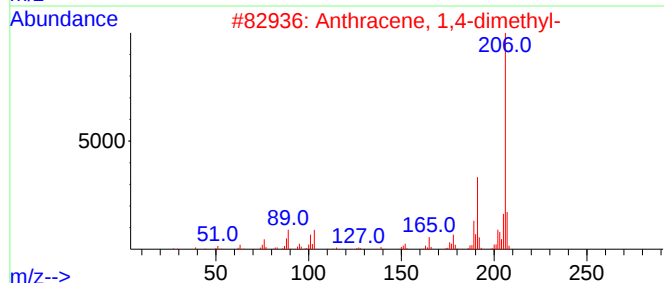
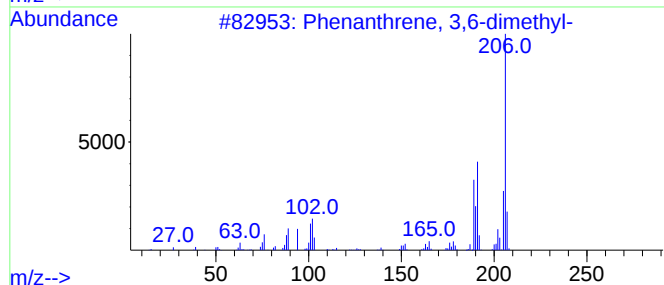
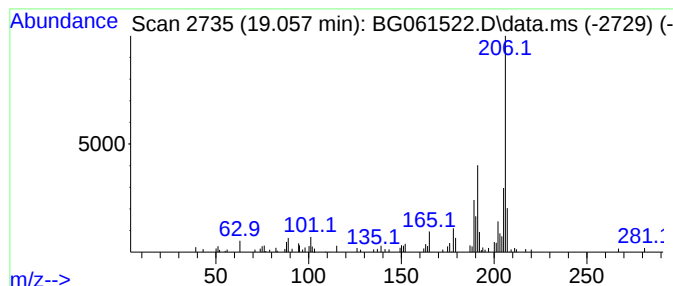
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Quant Title : SVOA CALIBRATION

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

Peak Number 11 Phenanthrene, 3,6-dimethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.057	4.69 ng/ul	106441	Phenanthrene-d10	17.259

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenanthrene, 3,6-dimethyl-	206	C16H14	001576-67-6	94
2			Anthracene, 1,4-dimethyl-	206	C16H14	000781-92-0	93
3			3,4-Dimethylphenanthrene	206	C16H14	1000452-24-5	93
4			Phenanthrene, 2,5-dimethyl-	206	C16H14	003674-66-6	93
5			di-p-Tolylacetylene	206	C16H14	002789-88-0	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG060624\
Data File : BG061522.D
Acq On : 6 Jun 2024 23:13
Operator : MA/JU
Sample : P2634-03
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BHBW9

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG052124.MA.M
Quant Title : SVOA CALIBRATION

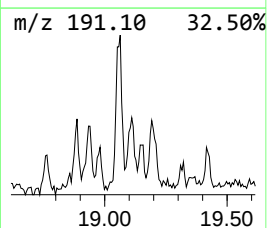
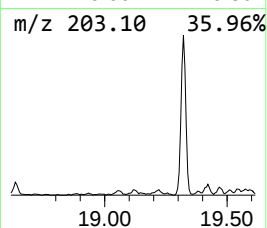
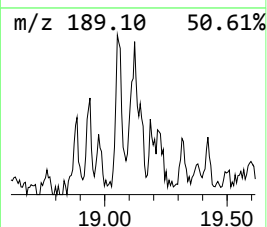
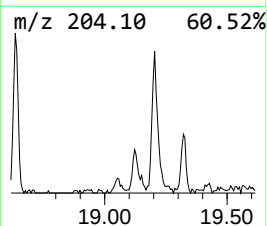
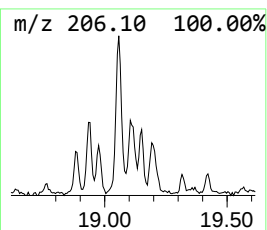
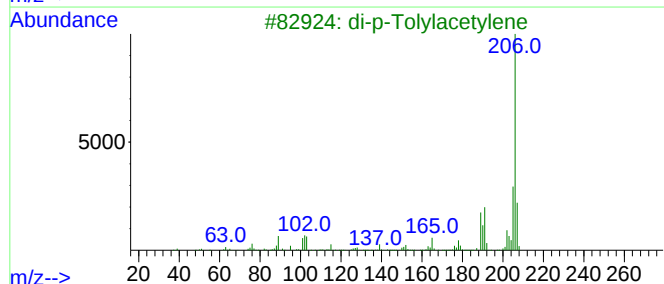
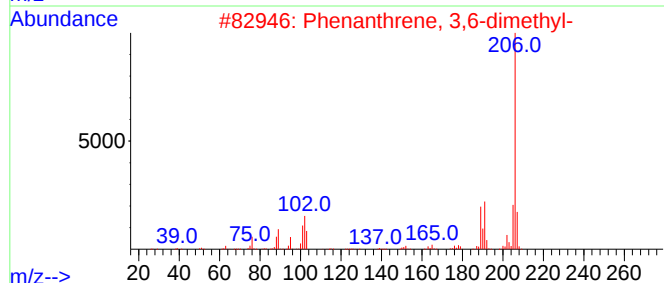
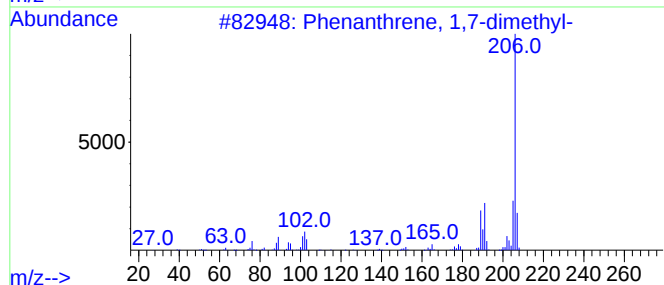
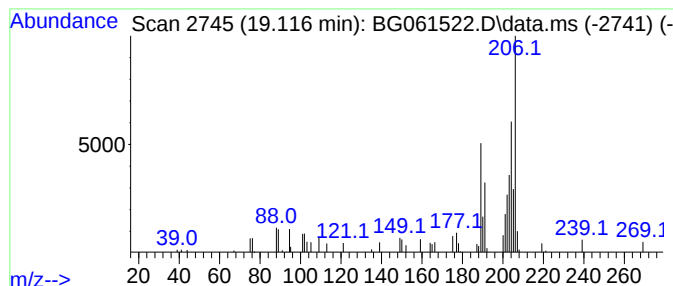
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 12 unknown-01

Concentration Rank 22

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.116	2.05 ng/ul	46421	Phenanthrene-d10	17.259

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenanthrene, 1,7-dimethyl-	206	C16H14	000483-87-4	49
2			Phenanthrene, 3,6-dimethyl-	206	C16H14	001576-67-6	46
3			di-p-Tolylacetylene	206	C16H14	002789-88-0	45
4			Anthracene, 1,4-dimethyl-	206	C16H14	000781-92-0	43
5			7-Methoxy-6-methyl-2,4-pteridine...	206	C8H10N6O	343347-40-0	41



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG060624\
Data File : BG061522.D
Acq On : 6 Jun 2024 23:13
Operator : MA/JU
Sample : P2634-03
Misc :
ALS Vial : 10 Sample Multiplier: 1

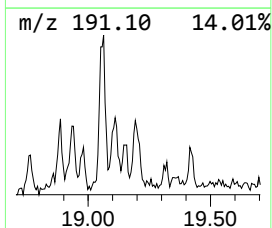
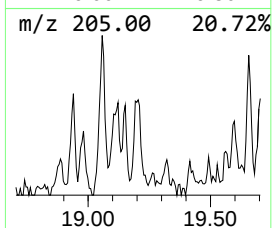
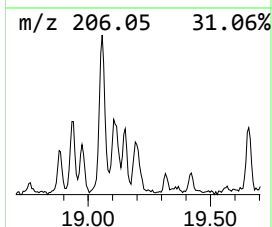
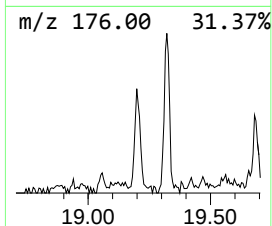
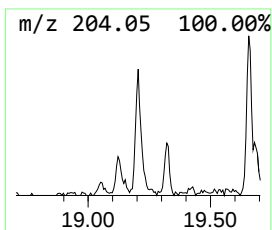
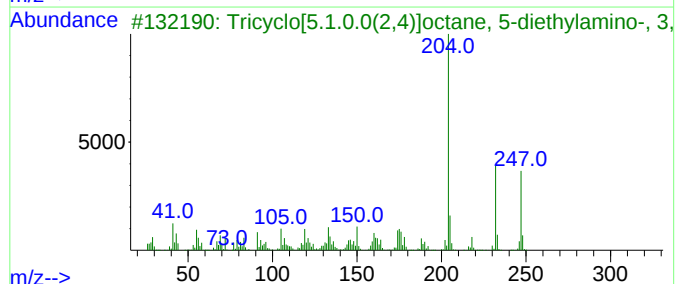
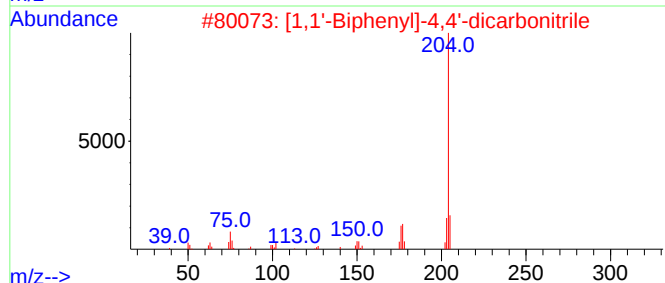
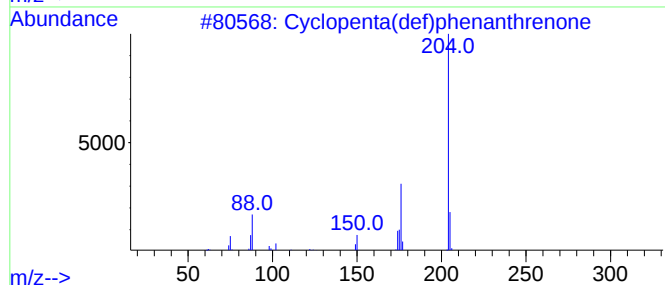
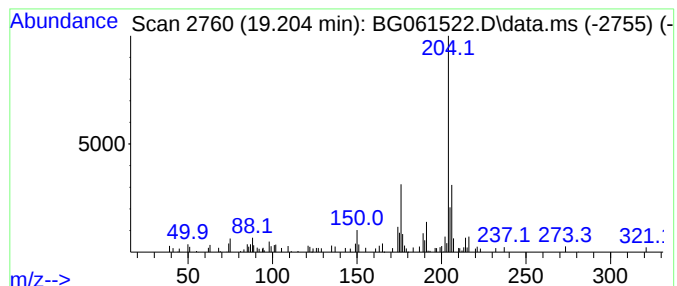
Instrument :
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG052124.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 13 Cyclopenta(def)phenanthrene Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.	
19.204	3.40 ng/ul	77156	Phenanthrene-d10	17.259	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclopenta(def)phenanthrenone	204	C15H8O	005737-13-3	76
2	[1,1'-Biphenyl]-4,4'-dicarbonitrile	204	C14H8N2	001591-30-6	52
3	Tricyclo[5.1.0.0(2,4)]octane, 5-...	247	C17H29N	1000063-59-3	50
4	[1,1'-Biphenyl]-2-ol, 5-chloro-	204	C12H9ClO	000607-12-5	50
5	3,4-Biphenyldicarbonitrile	204	C14H8N2	004128-63-6	50



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG060624\
Data File : BG061522.D
Acq On : 6 Jun 2024 23:13
Operator : MA/JU
Sample : P2634-03
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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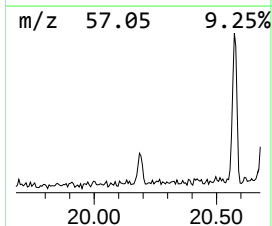
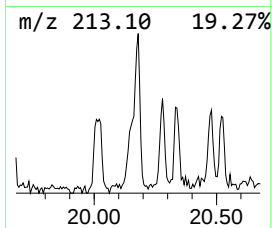
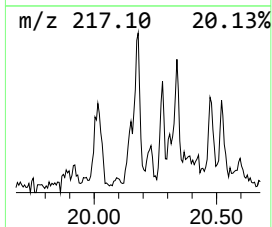
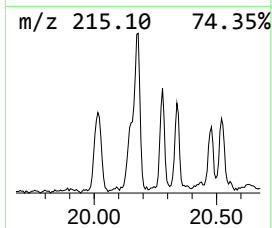
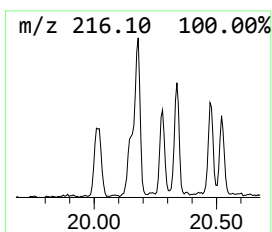
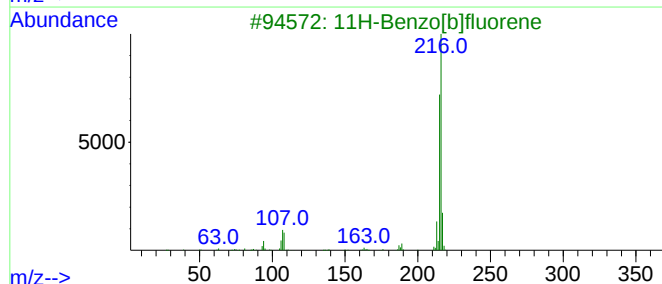
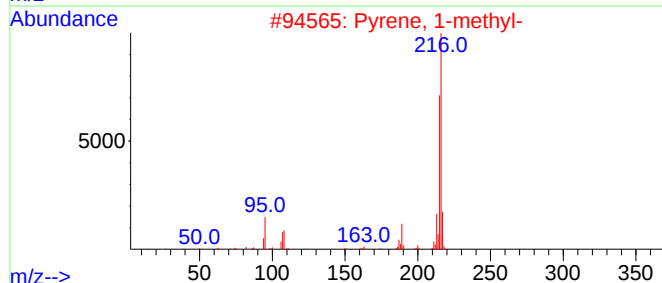
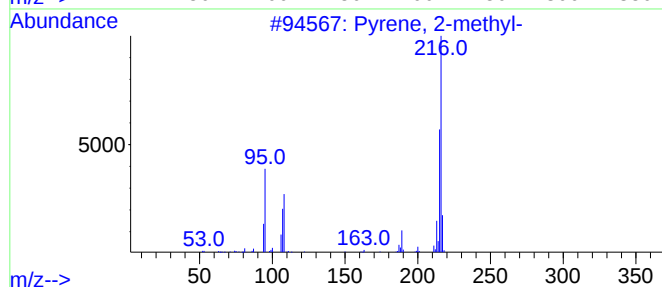
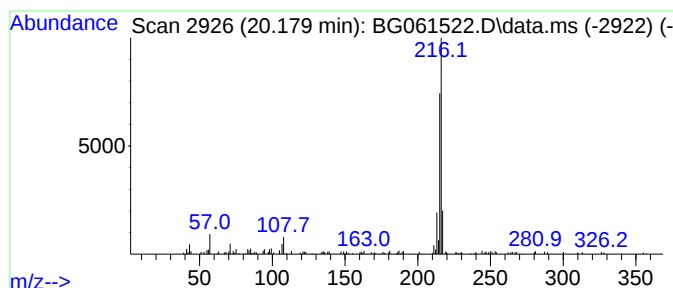
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Quant Title : SVOA CALIBRATION

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

Peak Number 14 Pyrene, 2-methyl- Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.179	2.70 ng/ul	159350	Chrysene-d12	21.519

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pyrene, 2-methyl-	216	C17H12	003442-78-2	91
2			Pyrene, 1-methyl-	216	C17H12	002381-21-7	91
3			11H-Benzo[b]fluorene	216	C17H12	000243-17-4	90
4			11H-Benzo[a]fluorene	216	C17H12	000238-84-6	90
5			Pyrene, 1-methyl-	216	C17H12	002381-21-7	86



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG060624\
Data File : BG061522.D
Acq On : 6 Jun 2024 23:13
Operator : MA/JU
Sample : P2634-03
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BHBW9

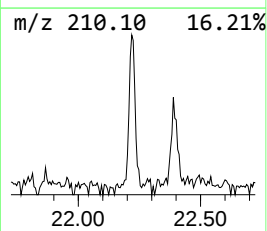
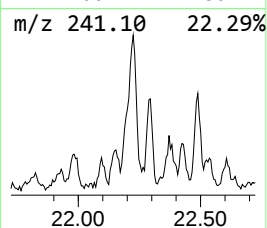
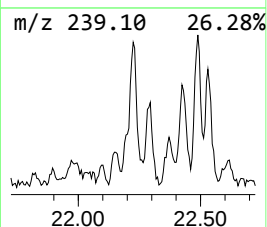
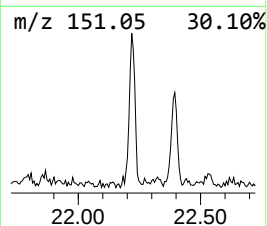
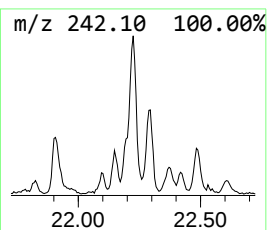
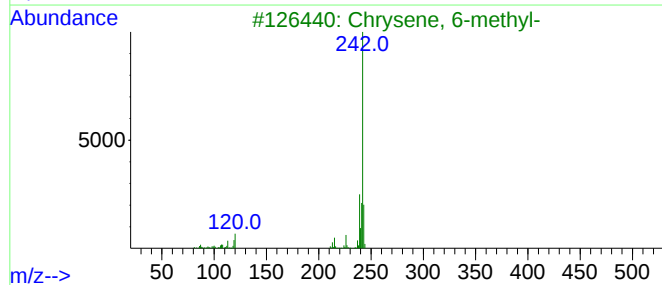
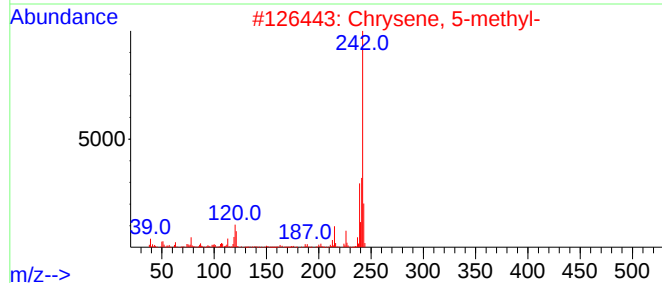
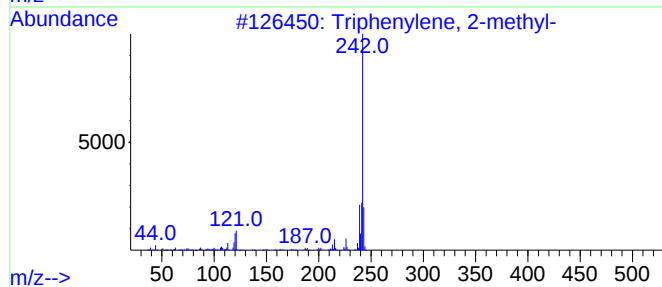
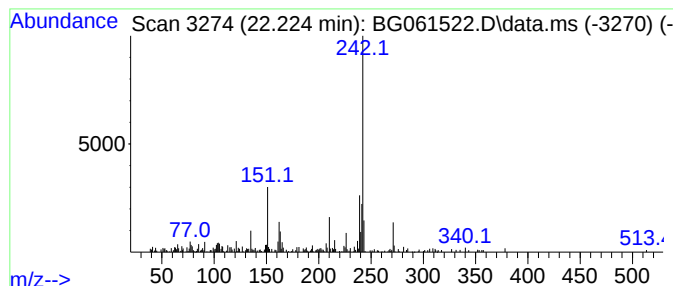
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Quant Title : SVOA CALIBRATION

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

Peak Number 15 Triphenylene, 2-methyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.224	3.22 ng/ul	189914	Chrysene-d12	21.519

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Triphenylene, 2-methyl-	242	C19H14	001705-84-6	83
2			Chrysene, 5-methyl-	242	C19H14	003697-24-3	64
3			Chrysene, 6-methyl-	242	C19H14	001705-85-7	60
4			Benz[a]anthracene, 2-methyl-	242	C19H14	002498-76-2	60
5			Benz[a]anthracene, 4-methyl-	242	C19H14	000316-49-4	60



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG060624\
Data File : BG061522.D
Acq On : 6 Jun 2024 23:13
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Sample : P2634-03
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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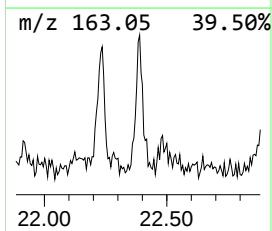
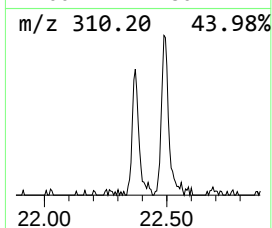
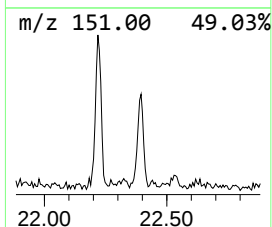
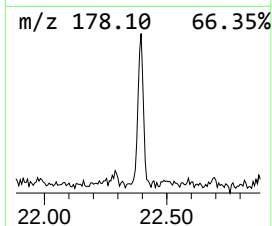
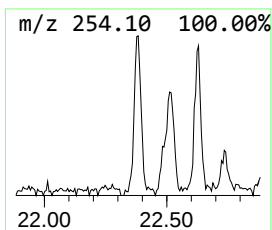
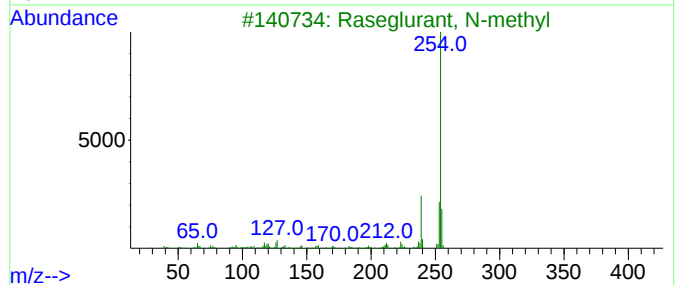
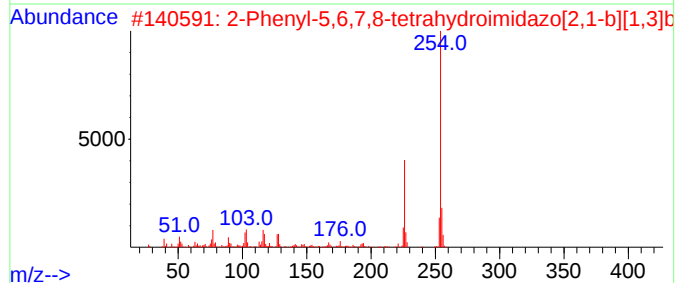
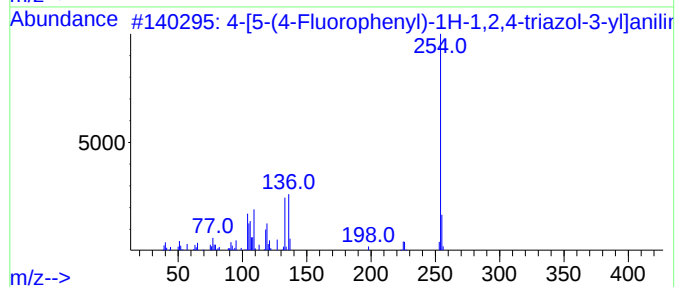
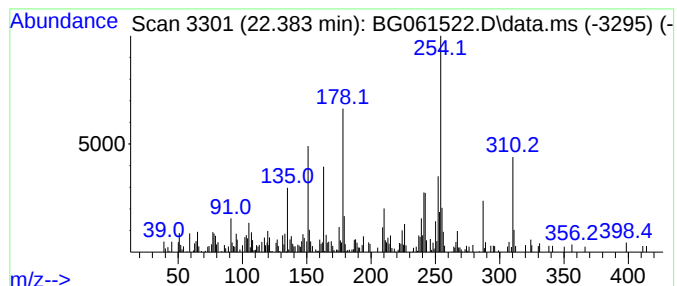
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Quant Title : SVOA CALIBRATION

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

Peak Number 16 unknown-02 Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.383	2.49 ng/ul	146826	Chrysene-d12	21.519

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	4-	[5-(4-Fluorophenyl)-1H-1,2,4-t...	254	C14H11FN4	1000387-34-8	35	
2	2-Phenyl-5,6,7,8-tetrahydroimida...	254	C15H14N2S	025752-17-4	35		
3	Raseglurant, N-methyl	254	C16H15FN2	1000498-91-9	35		
4	2(3H)-indolizine, 1-(4-nitroph...	254	C14H10N2O3	1000396-69-7	30		
5	trans-4'-(Methylthio)chalcone	254	C16H14OS	017129-07-6	30		



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Data File : BG061522.D
Acq On : 6 Jun 2024 23:13
Operator : MA/JU
Sample : P2634-03
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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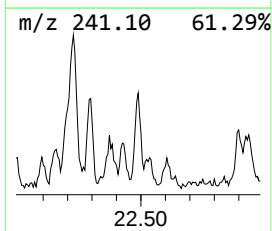
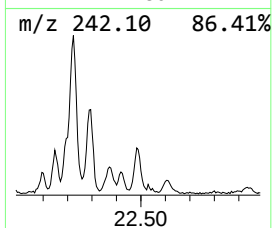
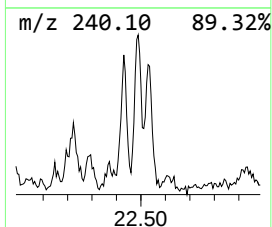
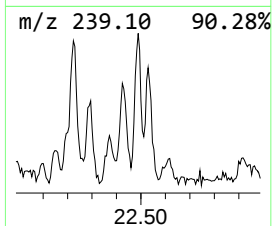
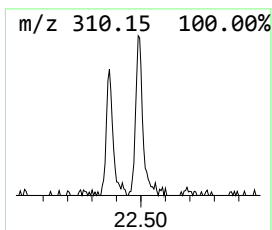
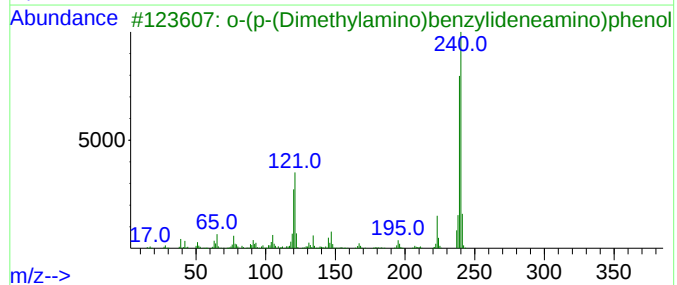
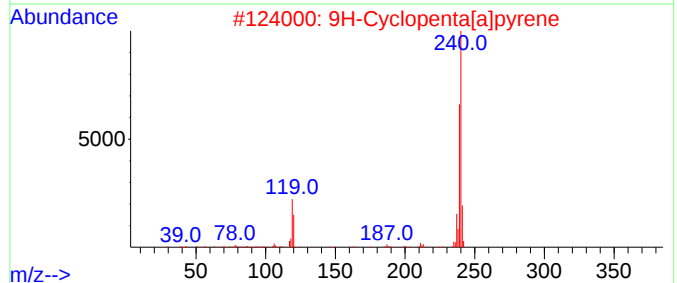
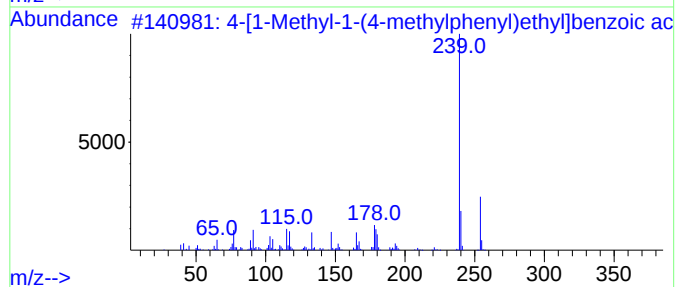
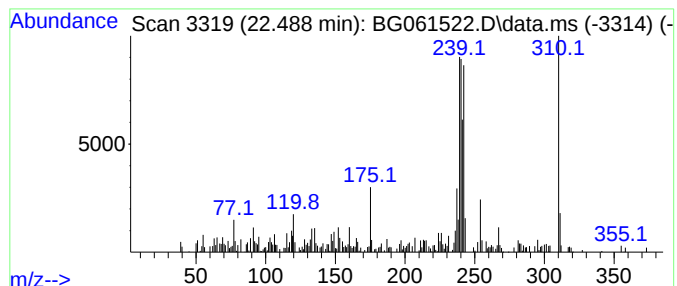
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Quant Title : SVOA CALIBRATION

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

Peak Number 17 unknown-03 Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.488	2.69 ng/ul	158667	Chrysene-d12	21.519

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	4-	[1-Methyl-1-(4-methylphenyl)et...	254	C17H18O2	006278-37-1	44	
2	9H-Cyclopenta[a]pyrene		240	C19H12	050861-05-7	41	
3	o-(p-(Dimethylamino)benzylidenea...		240	C15H16N2O	023837-35-6	38	
4	7-Bromo-4-methoxy-indan-1-one		240	C10H9BrO2	1010202-16-3	30	
5	1H-Pyrrolo[2,3-f]quinoline-8-car...		310	C19H22N2O2	199919-64-7	25	



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG060624\
Data File : BG061522.D
Acq On : 6 Jun 2024 23:13
Operator : MA/JU
Sample : P2634-03
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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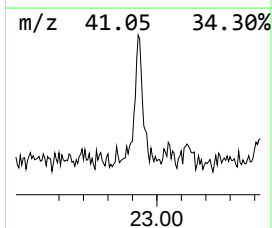
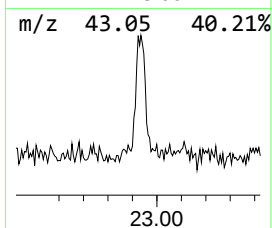
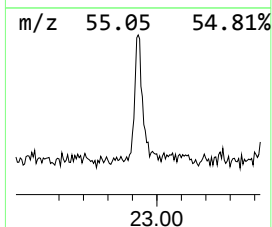
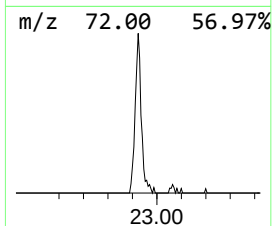
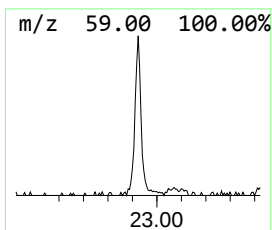
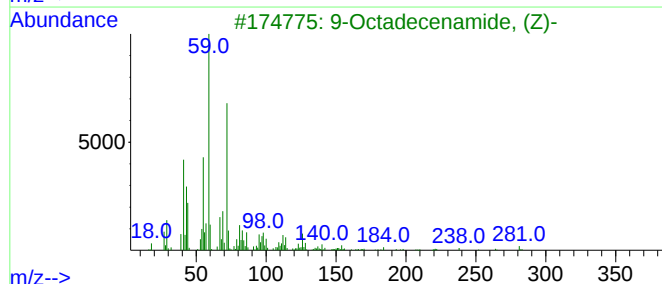
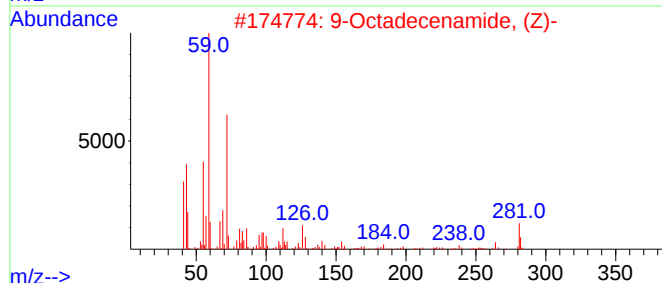
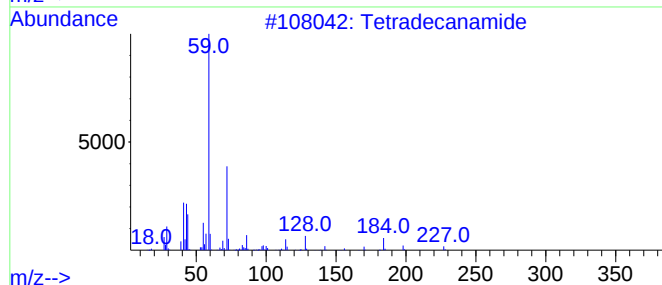
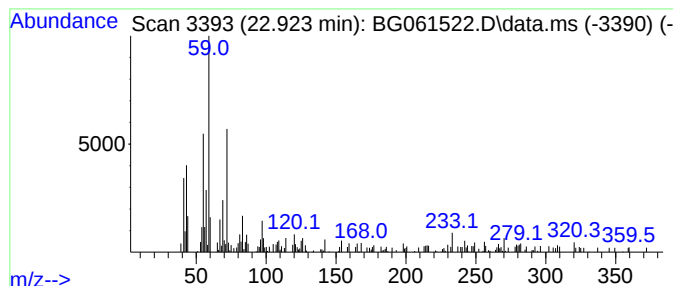
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Quant Title : SVOA CALIBRATION

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

Peak Number 18 Tetradecanamide Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.923	2.48 ng/ul	146185	Chrysene-d12	21.519

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tetradecanamide	227	C14H29NO	000638-58-4	70
2			9-Octadecenamide, (Z)-	281	C18H35NO	000301-02-0	68
3			9-Octadecenamide, (Z)-	281	C18H35NO	000301-02-0	68
4			13-Docosenamide, (Z)-	337	C22H43NO	000112-84-5	62
5			9-Octadecenamide, (Z)-	281	C18H35NO	000301-02-0	59



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG060624\
Data File : BG061522.D
Acq On : 6 Jun 2024 23:13
Operator : MA/JU
Sample : P2634-03
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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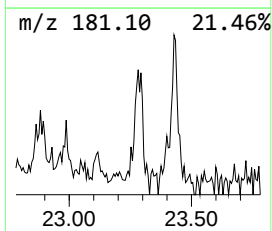
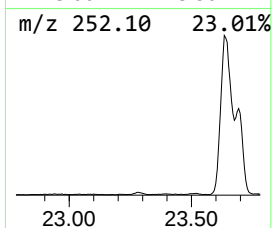
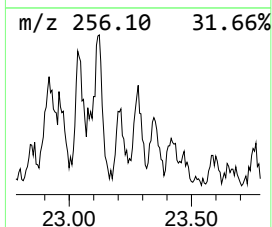
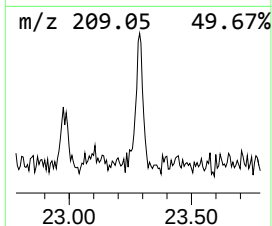
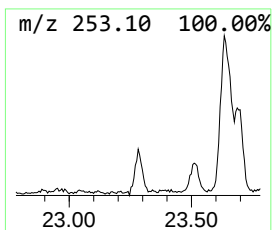
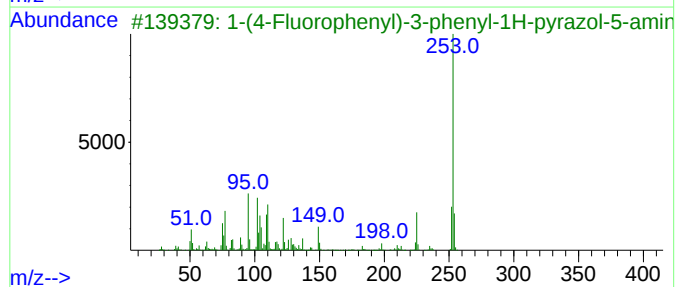
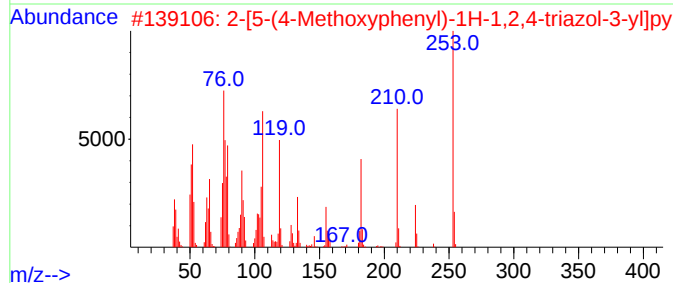
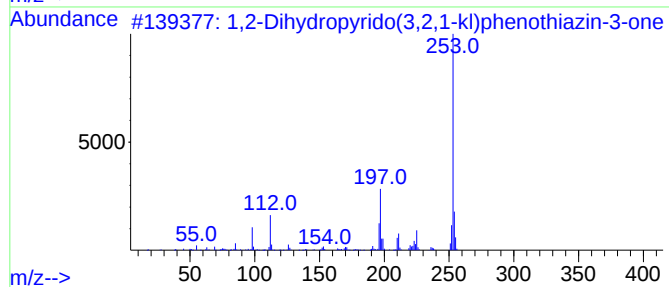
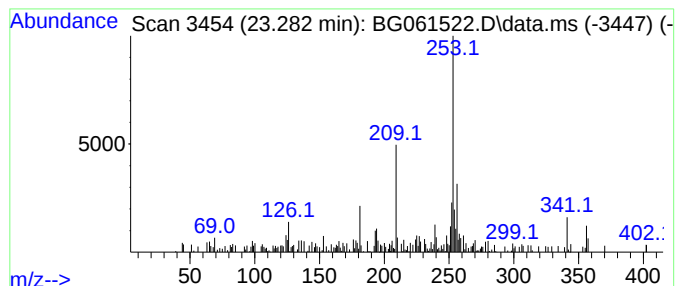
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Quant Title : SVOA CALIBRATION

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

Peak Number 19 unknown-04 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.282	4.33 ng/ul	117192	Perylene-d12	24.615

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1,2-Dihydropyrido(3,2,1-kl)pheno...	253	C15H11NOS	069513-42-4	43
2			2-[5-(4-Methoxyphenyl)-1H-1,2,4-...	253	C13H11N5O	159053-06-2	38
3			1-(4-Fluorophenyl)-3-phenyl-1H-p...	253	C15H12FN3	072411-51-9	38
4			Fumaric acid, 3-hexyl undecyl ester	354	C21H38O4	1000348-61-0	38
5			5,12-ethanonaphthacene-6-carboni...	281	C21H15N	1000397-19-0	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG060624\
Data File : BG061522.D
Acq On : 6 Jun 2024 23:13
Operator : MA/JU
Sample : P2634-03
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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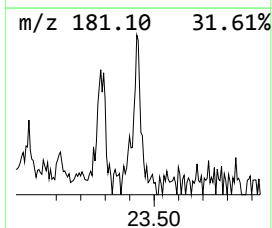
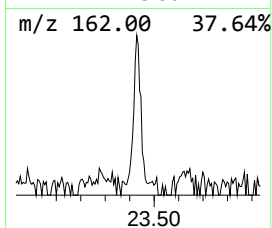
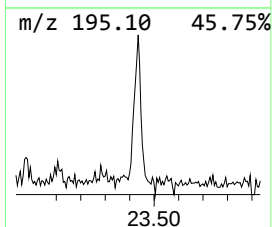
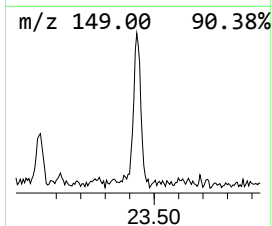
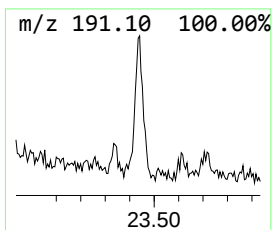
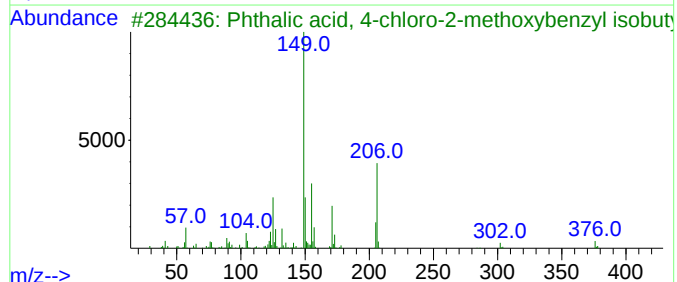
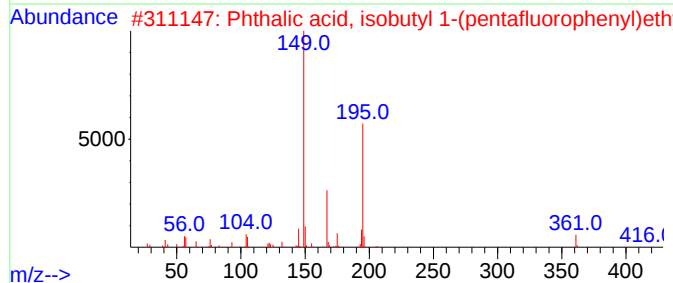
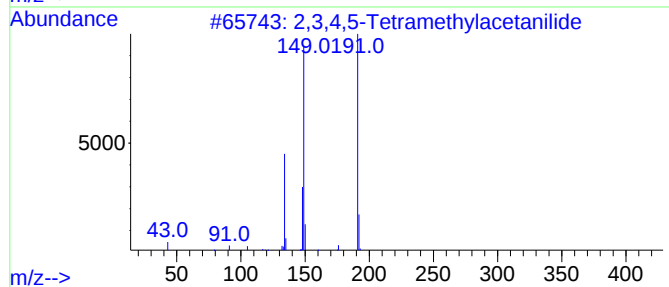
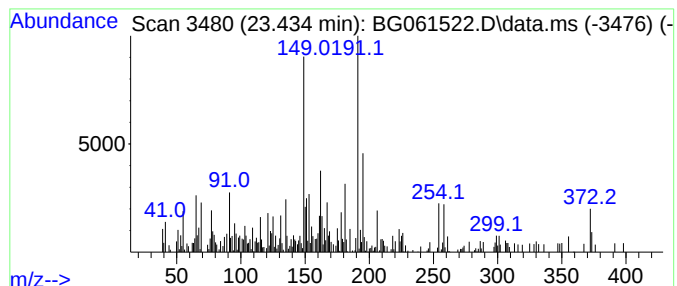
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Quant Title : SVOA CALIBRATION

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

Peak Number 20 unknown-05 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.434	7.90 ng/ul	213790	Perylene-d12	24.615

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,3,4,5-Tetramethylacetanilide	191	C12H17NO	1000432-93-7	43		
2	Phthalic acid, isobutyl 1-(penta...	416	C20H17F5O4	1000382-92-6	27		
3	Phthalic acid, 4-chloro-2-methox...	376	C20H21ClO5	1000382-83-7	25		
4	Phenol, 2,6-bis(1,1-dimethylethyl)-	206	C14H22O	000128-39-2	22		
5	Oxirane, [[4-(1,1-dimethylethyl)...	206	C13H18O2	003101-60-8	18		



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG060624\
Data File : BG061522.D
Acq On : 6 Jun 2024 23:13
Operator : MA/JU
Sample : P2634-03
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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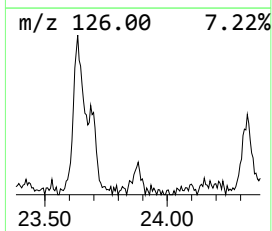
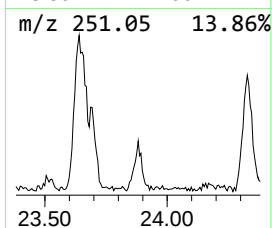
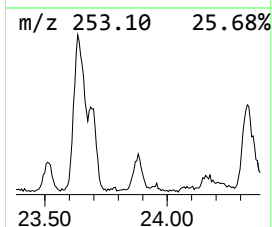
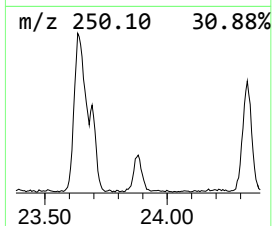
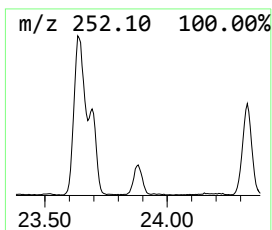
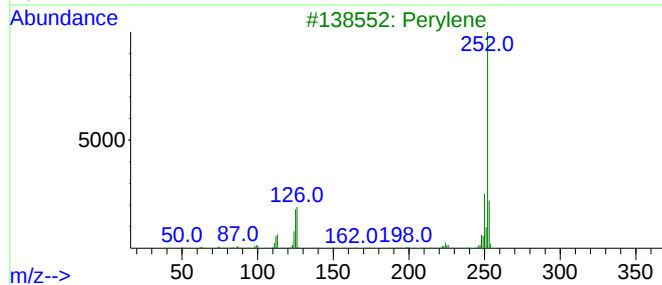
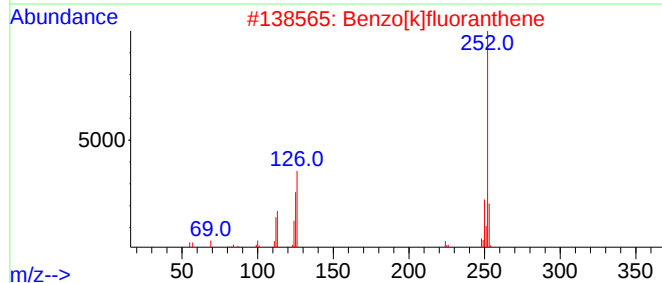
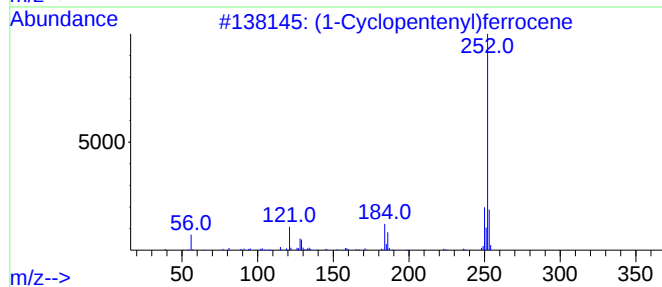
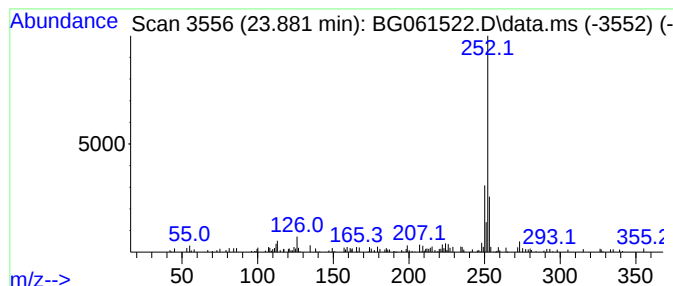
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Quant Title : SVOA CALIBRATION

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

Peak Number 21 (1-Cyclopentenyl)ferrocene Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.881	4.44 ng/ul	120325	Perylene-d12	24.615

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			(1-Cyclopentenyl)ferrocene	252	C15H16Fe	012260-67-2	80
2			Benzo[k]fluoranthene	252	C20H12	000207-08-9	76
3			Perylene	252	C20H12	000198-55-0	70
4			Benzo[a]pyrene	252	C20H12	000050-32-8	68
5			Benzo[b]fluoranthene	252	C20H12	000205-99-2	68



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG060624\
Data File : BG061522.D
Acq On : 6 Jun 2024 23:13
Operator : MA/JU
Sample : P2634-03
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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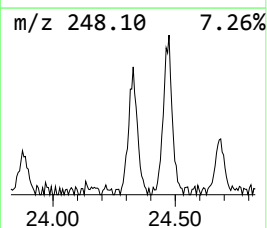
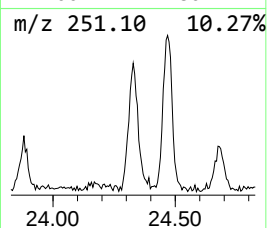
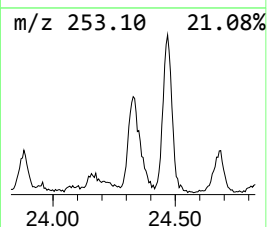
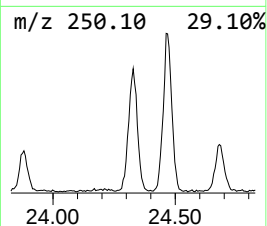
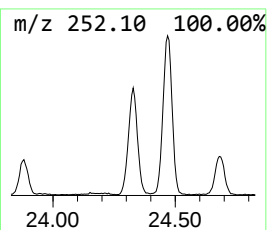
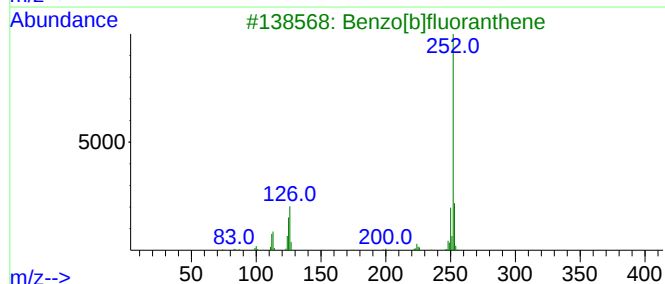
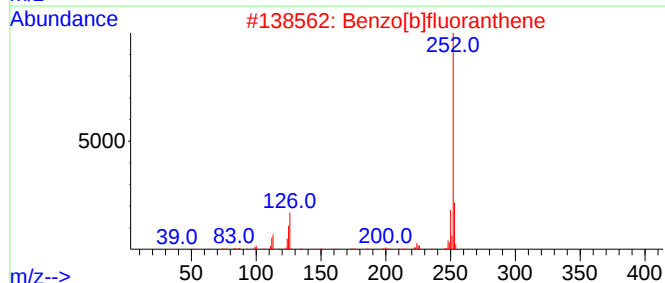
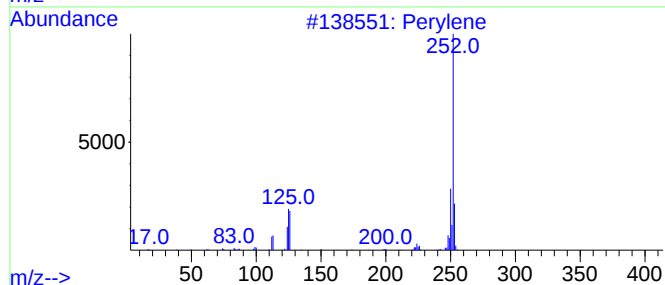
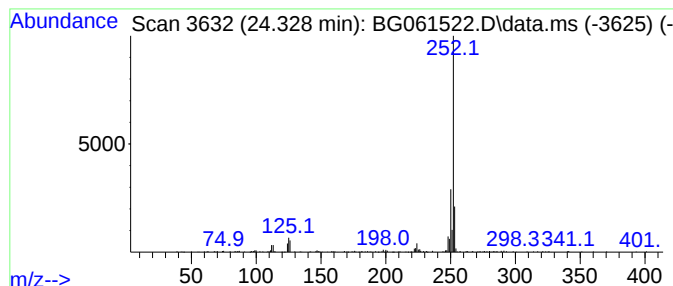
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Quant Title : SVOA CALIBRATION

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

Peak Number 22 Perylene Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.328	12.19 ng/ul	330029	Perylene-d12	24.615

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Perylene	252	C20H12	000198-55-0	89
2			Benzo[b]fluoranthene	252	C20H12	000205-99-2	81
3			Benzo[b]fluoranthene	252	C20H12	000205-99-2	76
4			Benzo[a]pyrene	252	C20H12	000050-32-8	74
5			Benzo[e]pyrene	252	C20H12	000192-97-2	74



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG060624\
 Data File : BG061522.D
 Acq On : 6 Jun 2024 23:13
 Operator : MA/JU
 Sample : P2634-03
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 BHBW9

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG052124.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
Butane, 2-metho...	3.111	37.3	ng/ul	405066	1	7.882	217025	20.0
5,5-Dimethyl-1-...	7.318	2.3	ng/ul	24499	1	7.882	217025	20.0
Pentanedioic ac...	9.586	2.0	ng/ul	27643	2	10.667	269446	20.0
1-Phenoxypropan...	11.413	2.5	ng/ul	33678	2	10.667	269446	20.0
Phenanthrene, 2...	18.135	5.2	ng/ul	116899	4	17.259	453541	20.0
Phenanthrene, 1...	18.188	5.6	ng/ul	127666	4	17.259	453541	20.0
4H-Cyclopenta[d...	18.329	5.7	ng/ul	129688	4	17.259	453541	20.0
Phenanthrene, 3...	18.370	2.7	ng/ul	60388	4	17.259	453541	20.0
9,10-di(Chlorom...	18.634	2.9	ng/ul	65360	4	17.259	453541	20.0
9,10-Anthracene...	18.687	3.5	ng/ul	78608	4	17.259	453541	20.0
Phenanthrene, 3...	19.057	4.7	ng/ul	106441	4	17.259	453541	20.0
unknown-01	19.116	2.0	ng/ul	46421	4	17.259	453541	20.0
Cyclopenta(def)...	19.204	3.4	ng/ul	77156	4	17.259	453541	20.0
Pyrene, 2-methyl-	20.179	2.7	ng/ul	159350	5	21.519	1180730	20.0
Triphenylene, 2...	22.224	3.2	ng/ul	189914	5	21.519	1180730	20.0
unknown-02	22.383	2.5	ng/ul	146826	5	21.519	1180730	20.0
unknown-03	22.488	2.7	ng/ul	158667	5	21.519	1180730	20.0
Tetradecanamide	22.923	2.5	ng/ul	146185	5	21.519	1180730	20.0
unknown-04	23.282	4.3	ng/ul	117192	6	24.615	541433	20.0
unknown-05	23.434	7.9	ng/ul	213790	6	24.615	541433	20.0
(1-Cyclopentenyl...	23.881	4.4	ng/ul	120325	6	24.615	541433	20.0
Perylene	24.328	12.2	ng/ul	330029	6	24.615	541433	20.0