

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM121523\
 Data File : BM043374.D
 Acq On : 16 Dec 2023 01:02
 Operator : MA/JU
 Sample : 05626-17
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 DCLG3

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_M\Methods\SFAM-EPA-BM121523.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BM043374.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.387	30	34	45	rVB	110135	150737	2.20%	0.123%
2	3.810	103	106	122	rBV	589311	952115	13.91%	0.777%
3	3.928	122	126	131	rVV	78402	110305	1.61%	0.090%
4	4.040	141	145	152	rVB	52671	83428	1.22%	0.068%
5	4.140	158	162	166	rVB	37341	48168	0.70%	0.039%
6	4.710	255	259	264	rBV4	18913	28417	0.42%	0.023%
7	5.069	316	320	328	rVB	63336	101487	1.48%	0.083%
8	5.181	331	339	344	rBV5	14612	27711	0.40%	0.023%
9	6.634	583	586	597	rVB5	19724	38796	0.57%	0.032%
10	7.151	666	674	685	rBV	1805642	2939095	42.94%	2.398%
11	7.334	698	705	712	rBV	1367204	2164810	31.63%	1.766%
12	7.398	712	716	724	rVB	76837	122391	1.79%	0.100%
13	7.522	730	737	744	rBV3	1931208	3431040	50.13%	2.799%
14	7.587	744	748	756	rVV	485286	722633	10.56%	0.590%
15	7.775	774	780	793	rBV	1541537	2528354	36.94%	2.063%
16	7.992	810	817	825	rBV	1703560	2778682	40.60%	2.267%
17	8.534	903	909	919	rBV	36317	75729	1.11%	0.062%
18	8.704	929	938	947	rBV	2251789	3567293	52.12%	2.910%
19	8.828	953	959	966	rVB	179787	289784	4.23%	0.236%
20	9.163	1008	1016	1028	rBV	1628702	2754546	40.24%	2.247%
21	9.263	1028	1033	1040	rVV	12548	28549	0.42%	0.023%
22	9.551	1077	1082	1092	rVB4	16416	39758	0.58%	0.032%
23	9.739	1109	1114	1120	rVB2	37424	63327	0.93%	0.052%
24	9.886	1128	1139	1148	rBV	1346964	2256452	32.97%	1.841%
25	10.022	1155	1162	1169	rBV3	31146	64757	0.95%	0.053%
26	10.422	1222	1230	1239	rBV	2084919	3463675	50.60%	2.826%
27	10.804	1285	1295	1301	rBV	2318457	3965324	57.93%	3.235%
28	10.851	1301	1303	1309	rVB	68888	97186	1.42%	0.079%
29	10.945	1311	1319	1334	rBV	1559699	2665521	38.94%	2.175%
30	11.345	1380	1387	1394	rBV10	22786	51925	0.76%	0.042%
31	11.551	1415	1422	1426	rBV2	31154	59312	0.87%	0.048%
32	11.992	1489	1497	1503	rBV6	11958	34033	0.50%	0.028%
33	12.216	1530	1535	1539	rVB5	22823	36951	0.54%	0.030%
34	12.304	1545	1550	1558	rBV2	36579	78108	1.14%	0.064%
35	12.386	1558	1564	1571	rVB	39047	72754	1.06%	0.059%
36	12.457	1571	1576	1581	rBV	36402	57640	0.84%	0.047%

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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_M\Methods\SFAM-EPA-BM121523.MA.M
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37	12.586	1592	1598	1605	rBV8	17630	40512	0.59%	0.033%
38	12.675	1608	1613	1619	rVB2	20795	37220	0.54%	0.030%
39	13.451	1740	1745	1751	rVV2	42480	78564	1.15%	0.064%
40	13.522	1751	1757	1767	rVV	96921	213604	3.12%	0.174%
41	13.945	1824	1829	1834	rVB6	35400	63872	0.93%	0.052%
42	14.039	1834	1845	1854	rBV	3107732	4507000	65.85%	3.677%
43	14.169	1862	1867	1874	rVB10	17019	41547	0.61%	0.034%
44	14.316	1885	1892	1907	rVB	3805854	5895226	86.13%	4.810%
45	14.621	1935	1944	1951	rBV	3168951	4643309	67.84%	3.788%
46	14.686	1951	1955	1959	rVB	50471	68001	0.99%	0.055%
47	14.769	1965	1969	1972	rBV6	20806	30152	0.44%	0.025%
48	14.821	1972	1978	1989	rBV	1349313	2051990	29.98%	1.674%
49	14.980	2001	2005	2008	rBV3	44183	59189	0.86%	0.048%
50	15.021	2008	2012	2015	rBV	48481	78863	1.15%	0.064%
51	15.163	2031	2036	2040	rVB8	14978	27184	0.40%	0.022%
52	15.410	2074	2078	2082	rBV5	14010	29922	0.44%	0.024%
53	15.616	2106	2113	2119	rBV	4510798	6456386	94.33%	5.268%
54	15.668	2119	2122	2127	rVB3	73205	100406	1.47%	0.082%
55	15.733	2127	2133	2140	rBV	1240465	1668145	24.37%	1.361%
56	16.351	2228	2238	2243	rVV4	81881	191786	2.80%	0.156%
57	16.668	2281	2292	2297	rBV4	24263	69310	1.01%	0.057%
58	17.010	2345	2350	2357	rVB3	86305	156122	2.28%	0.127%
59	17.121	2367	2369	2372	rVV3	20866	28179	0.41%	0.023%
60	17.180	2375	2379	2386	rVB4	37446	64572	0.94%	0.053%
61	17.286	2394	2397	2401	rVB3	24367	29603	0.43%	0.024%
62	17.362	2404	2410	2415	rBV	3076285	4459765	65.16%	3.639%
63	17.410	2415	2418	2422	rVB	283801	357419	5.22%	0.292%
64	17.462	2422	2427	2432	rBV	4191232	6084723	88.90%	4.964%
65	17.557	2438	2443	2452	rVB2	74410	141572	2.07%	0.116%
66	17.768	2472	2479	2492	rVB	175328	340180	4.97%	0.278%
67	18.262	2557	2563	2574	rBV2	357065	628481	9.18%	0.513%
68	18.368	2578	2581	2586	rVB	66033	97244	1.42%	0.079%
69	18.433	2587	2592	2598	rBV2	284059	488803	7.14%	0.399%
70	18.557	2609	2613	2616	rBV5	30262	48031	0.70%	0.039%
71	18.692	2634	2636	2641	rVB3	46766	52322	0.76%	0.043%
72	18.745	2641	2645	2648	rVV	61616	72301	1.06%	0.059%
73	18.780	2648	2651	2656	rVB	110784	149178	2.18%	0.122%
74	18.980	2681	2685	2690	rVB	342090	442083	6.46%	0.361%
75	19.115	2705	2708	2711	rBV	51106	69733	1.02%	0.057%
76	19.156	2711	2715	2719	rVV2	108921	167462	2.45%	0.137%

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Integration Parameters: LSCINT.P
 Integrator: RTE
 Smoothing : OFF Filtering: 5
 Sampling : 1 Min Area: 0.1 % of largest Peak
 Start Thrs: 0.2 Max Peaks: 100
 Stop Thrs : 0 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
 Peak separation: 5

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77	19.298	2735	2739	2740	rBV	91722	117058	1.71%	0.096%
78	19.415	2754	2759	2764	rVB	2432653	3229460	47.18%	2.635%
79	19.474	2764	2769	2774	rBV	707575	881494	12.88%	0.719%
80	19.751	2810	2816	2819	rBV	4723553	6844532	100.00%	5.584%
81	19.780	2819	2821	2829	rVV	2630496	2899832	42.37%	2.366%
82	19.851	2829	2833	2839	rVV2	104038	172850	2.53%	0.141%
83	20.103	2871	2876	2882	rBV4	193272	433632	6.34%	0.354%
84	20.268	2894	2904	2908	rBV3	325429	907124	13.25%	0.740%
85	20.368	2918	2921	2925	rBV	257652	330707	4.83%	0.270%
86	20.427	2925	2931	2935	rVV2	280025	575758	8.41%	0.470%
87	20.568	2952	2955	2958	rBV	211058	237318	3.47%	0.194%
88	21.015	3028	3031	3038	rVB2	263000	394757	5.77%	0.322%
89	21.186	3057	3060	3063	rVB	238843	262820	3.84%	0.214%
90	21.221	3063	3066	3069	rBV2	272465	306399	4.48%	0.250%
91	21.256	3069	3072	3078	rVB2	753561	1006744	14.71%	0.821%
92	21.539	3113	3120	3124	rBV2	3143801	5761809	84.18%	4.701%
93	21.574	3124	3126	3134	rVB	1454055	1918557	28.03%	1.565%
94	22.180	3226	3229	3235	rVB4	277095	447191	6.53%	0.365%
95	23.221	3400	3406	3413	rBV	2209844	5550617	81.10%	4.529%
96	23.744	3490	3495	3500	rBV	979281	1997082	29.18%	1.629%
97	23.803	3500	3505	3511	rVV2	2559148	5081679	74.24%	4.146%
98	23.856	3511	3514	3523	rVB2	649292	1089307	15.91%	0.889%
99	23.962	3525	3532	3538	rBV	1708015	3584753	52.37%	2.925%
100	26.485	3953	3961	3979	rVB2	606422	2357631	34.45%	1.923%

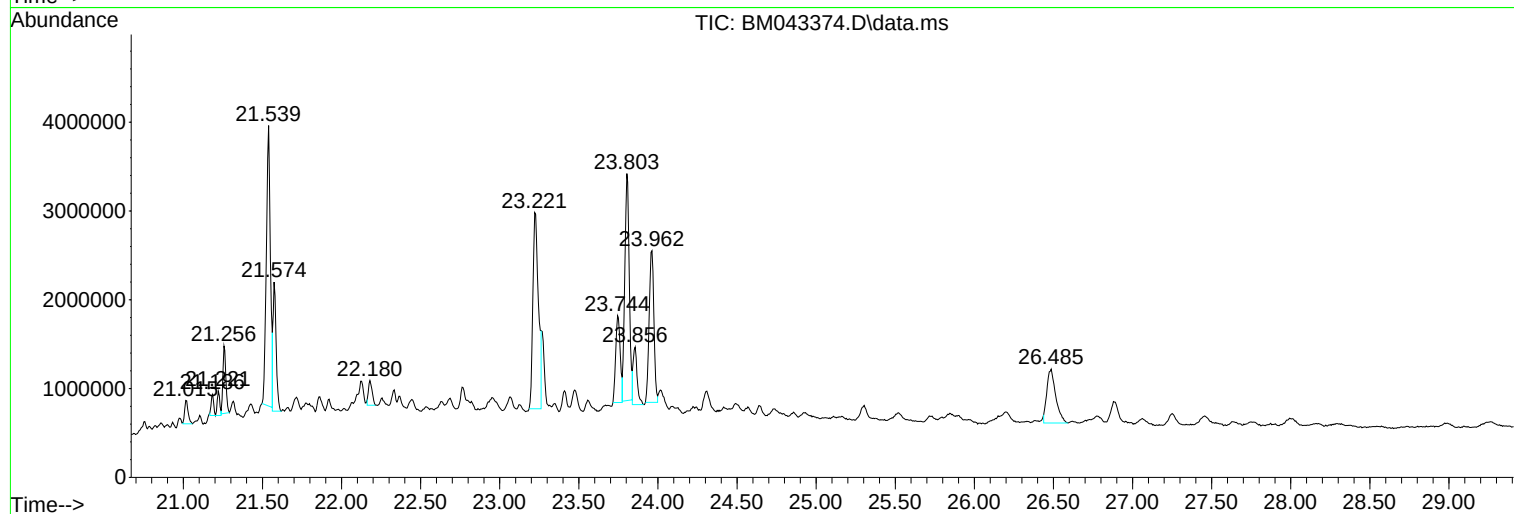
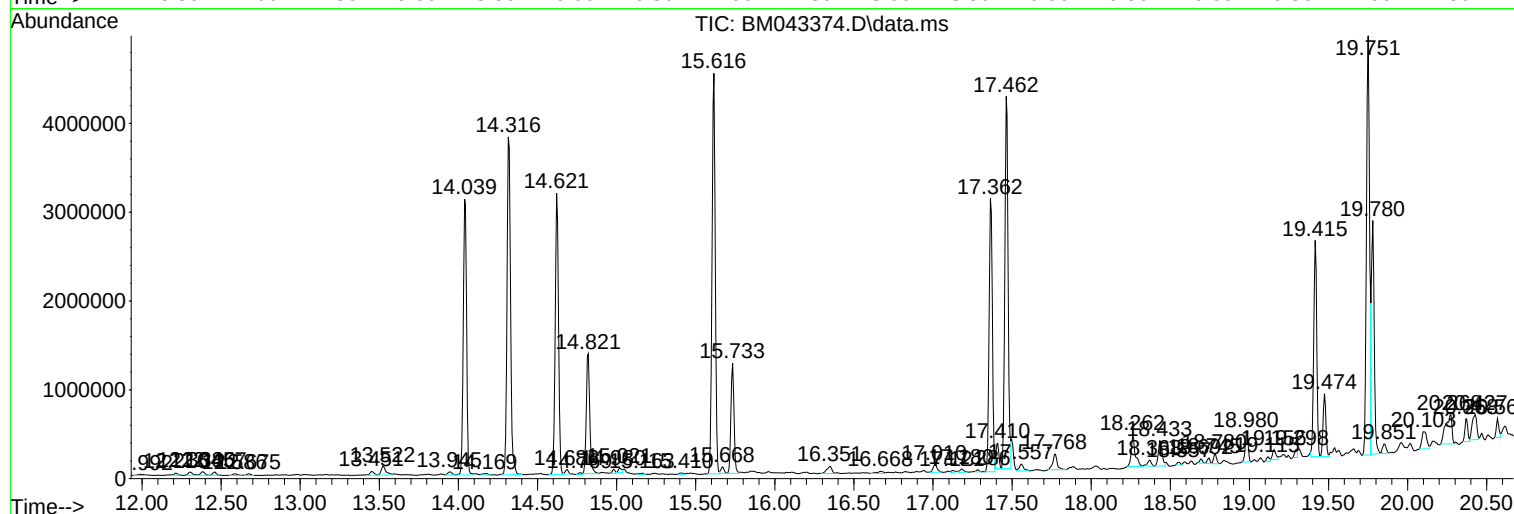
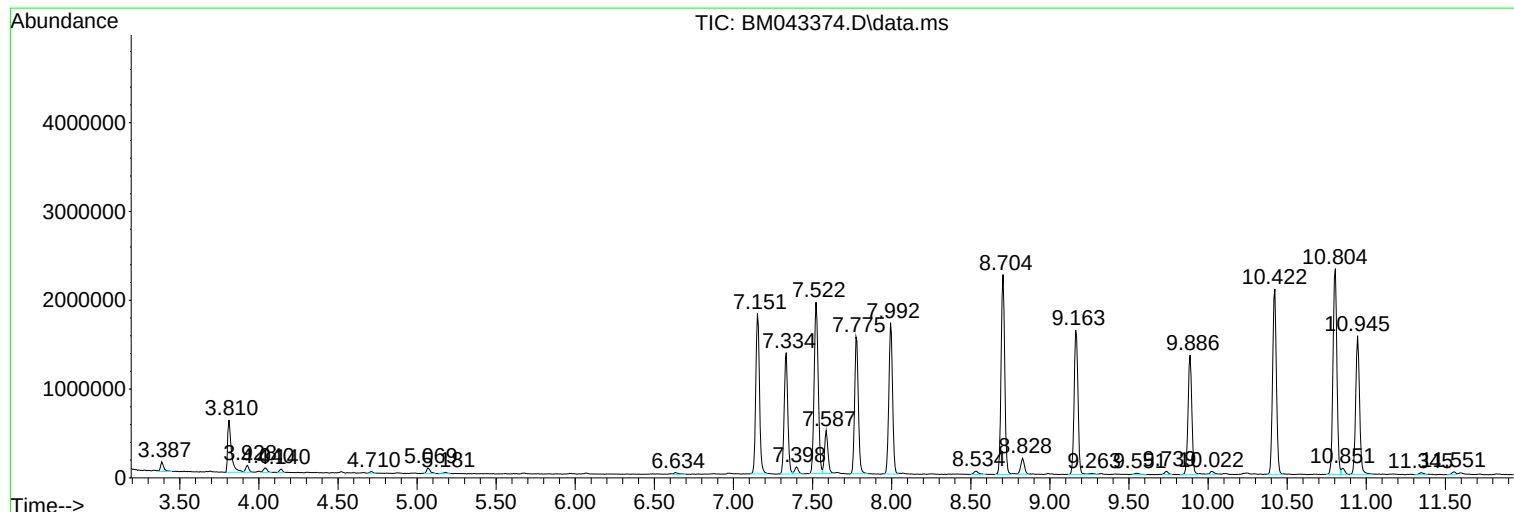
Sum of corrected areas: 122569865

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM121523.MA.M
Quant Title : SVOA CALIBRATION

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



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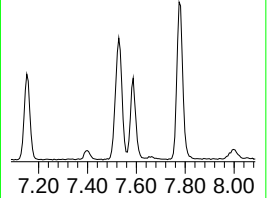
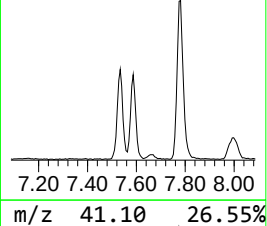
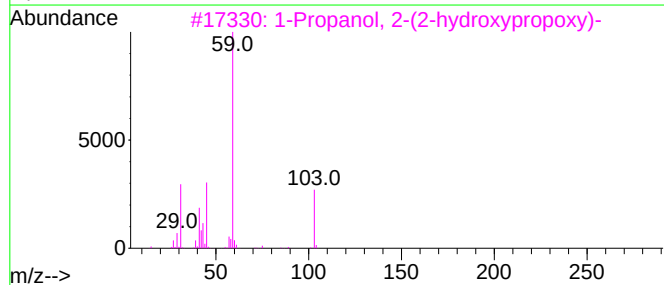
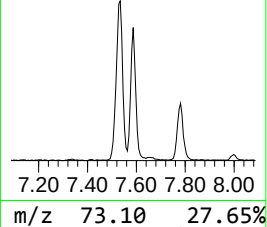
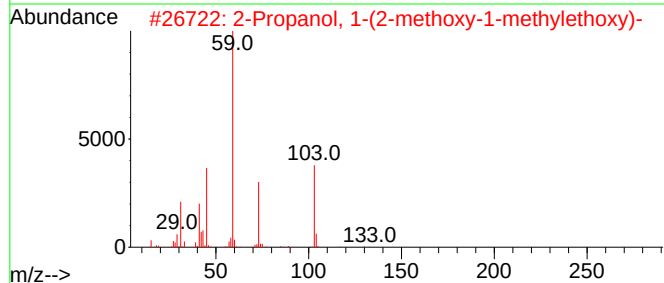
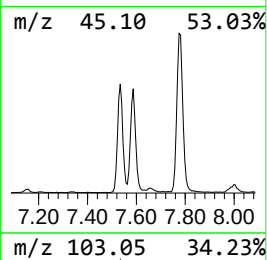
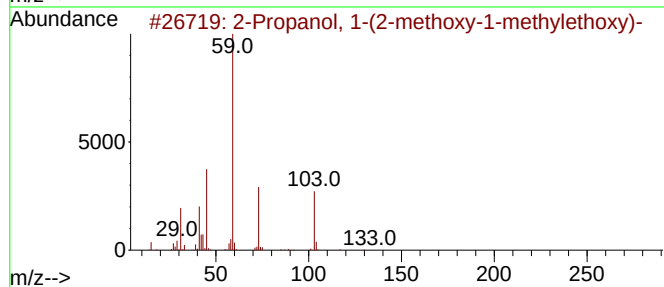
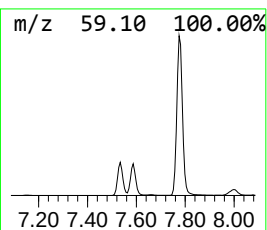
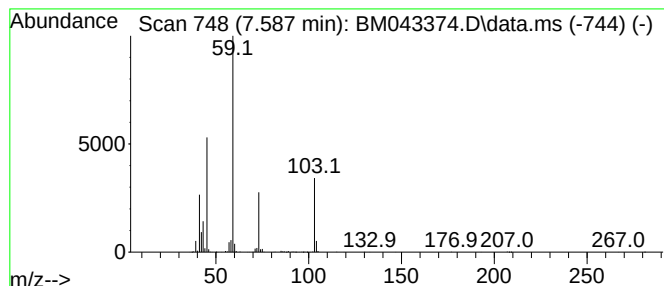
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM121523.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 2-Propanol, 1-(2-methoxy-1-... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.587	5.20 ng/ul	722633	1,4-Dichlorobenzene-d4	7.992

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Propanol, 1-(2-methoxy-1-methy...	148	C7H16O3	020324-32-7	83		
2	2-Propanol, 1-(2-methoxy-1-methy...	148	C7H16O3	020324-32-7	78		
3	1-Propanol, 2-(2-hydroxypropoxy)-	134	C6H14O3	000106-62-7	59		
4	1-Propanol, 2-(2-hydroxypropoxy)-	134	C6H14O3	000106-62-7	59		
5	1-Propanol, 2,2'-oxybis-	134	C6H14O3	000108-61-2	59		



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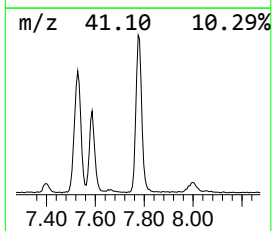
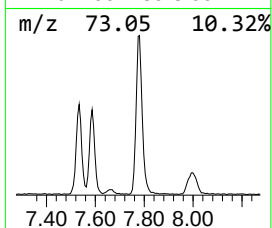
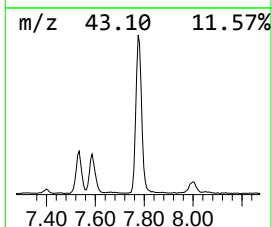
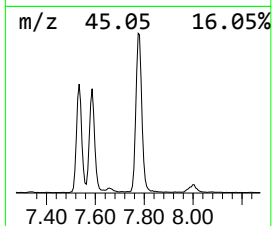
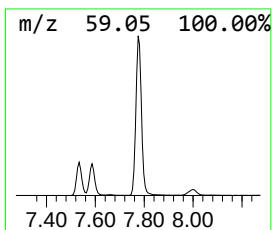
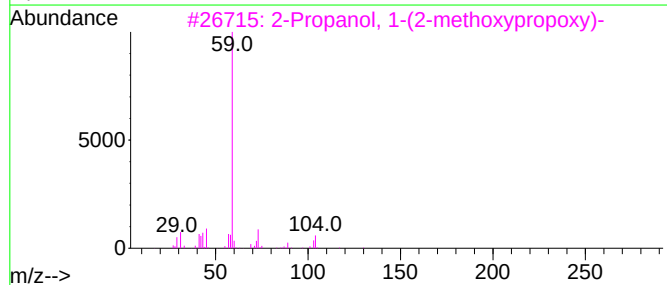
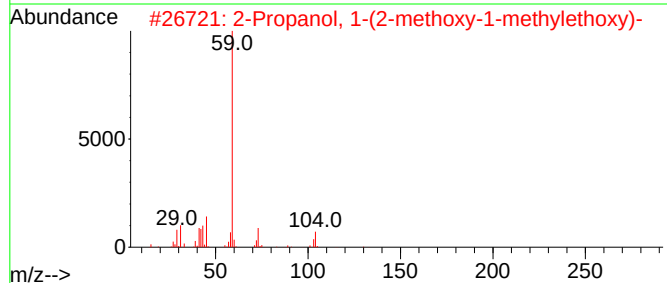
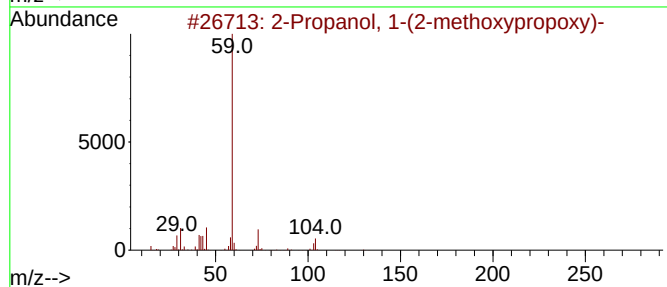
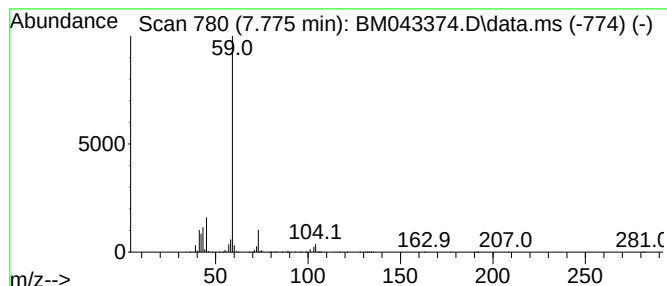
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM121523.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 2-Propanol, 1-(2-methoxypro... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.775	18.20 ng/ul	2528350	1,4-Dichlorobenzene-d4	7.992

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Propanol, 1-(2-methoxypropoxy)-	148	C7H16O3	013429-07-7	83	
2	2-Propanol, 1-(2-methoxy-1-methy...	148	C7H16O3	020324-32-7	74	
3	2-Propanol, 1-(2-methoxypropoxy)-	148	C7H16O3	013429-07-7	72	
4	1-Propanol, 2,2'-oxybis-	134	C6H14O3	000108-61-2	64	
5	1-Propanol, 2-(2-hydroxypropoxy)-	134	C6H14O3	000106-62-7	64	



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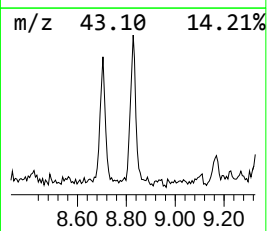
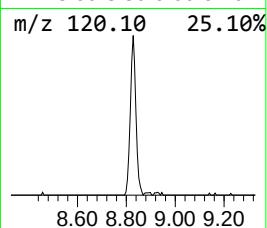
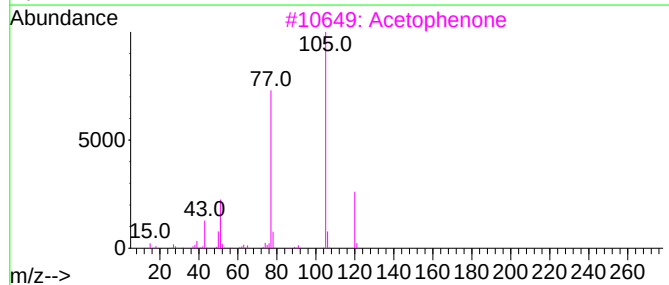
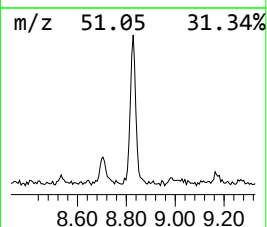
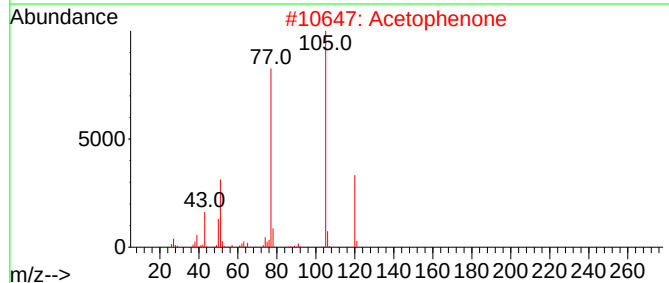
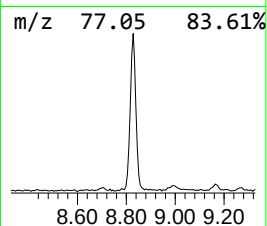
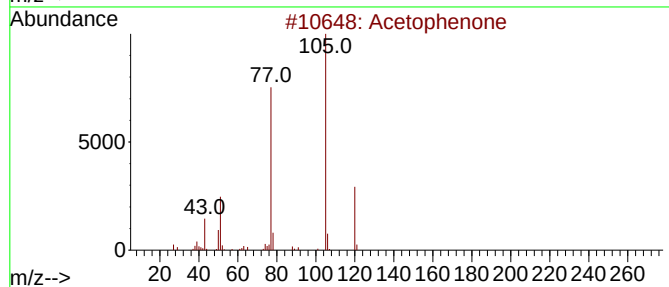
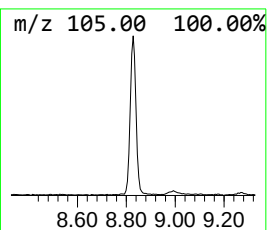
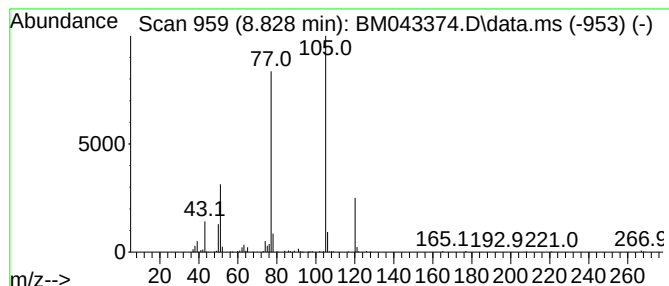
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Quant Title : SVOA CALIBRATION

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

Peak Number 3 Aceto phenone Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.828	2.09 ng/ul	289784	1,4-Dichlorobenzene-d4	7.992

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Acetophenone			120	C8H8O	000098-86-2	94
2	Acetophenone			120	C8H8O	000098-86-2	91
3	Acetophenone			120	C8H8O	000098-86-2	91
4	Acetophenone			120	C8H8O	000098-86-2	91
5	Acetophenone			120	C8H8O	000098-86-2	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM121523\
Data File : BM043374.D
Acq On : 16 Dec 2023 01:02
Operator : MA/JU
Sample : 05626-17
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
DCLG3

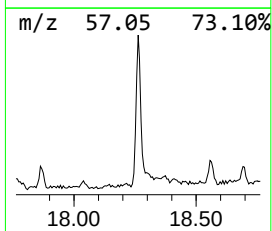
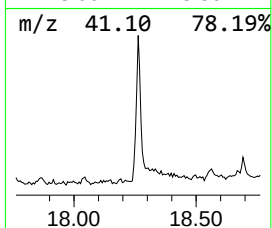
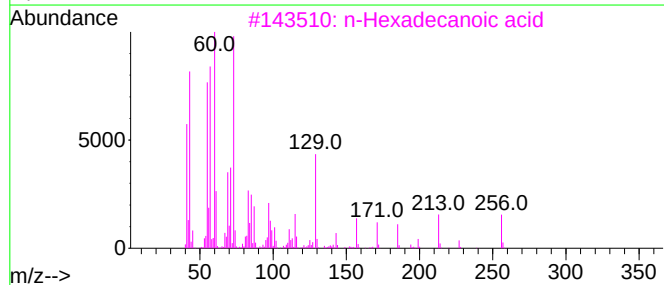
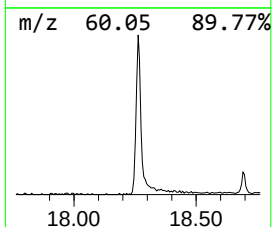
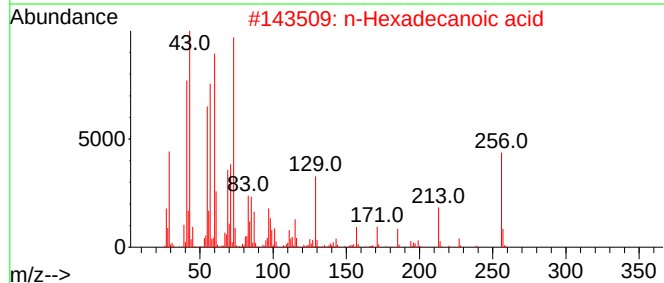
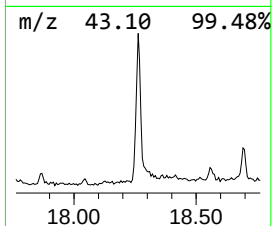
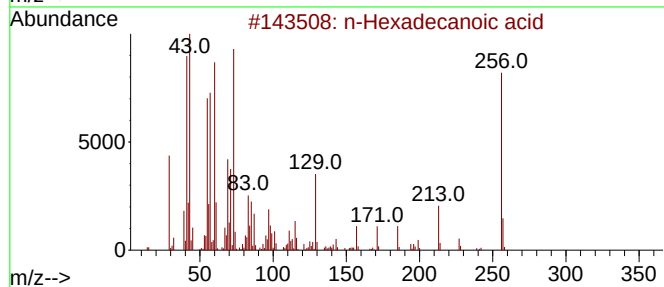
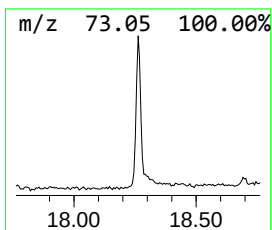
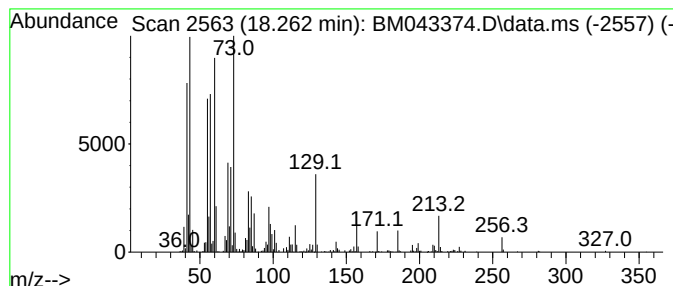
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 n-Hexadecanoic acid Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.262	2.82 ng/ul	628481	Phenanthrene-d10	17.363

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
3			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
4			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
5			Tetradecanoic acid	228	C14H28O2	000544-63-8	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM121523\
Data File : BM043374.D
Acq On : 16 Dec 2023 01:02
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Sample : 05626-17
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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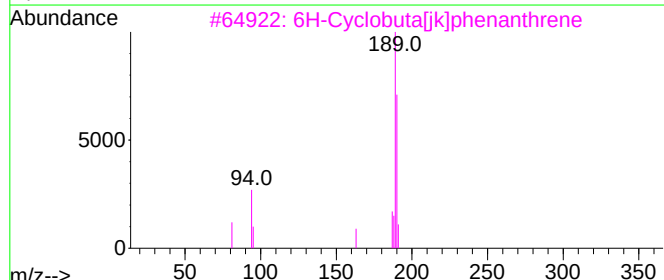
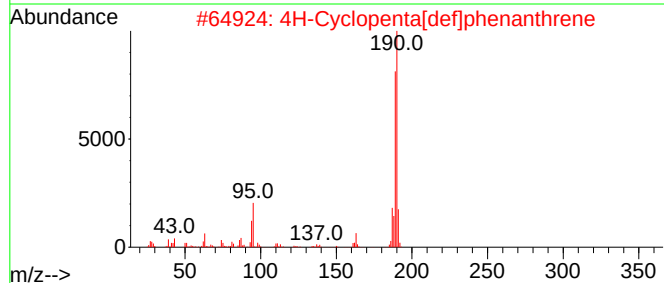
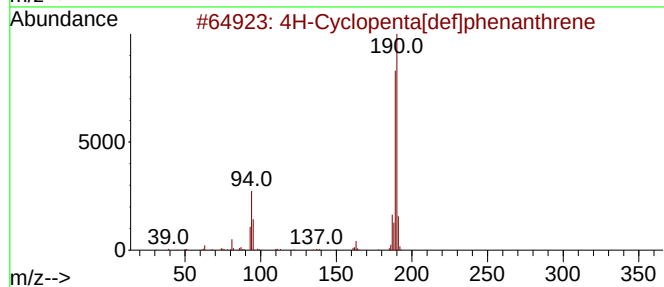
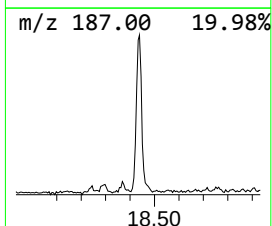
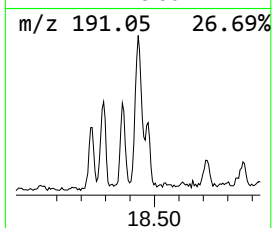
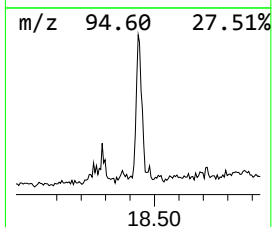
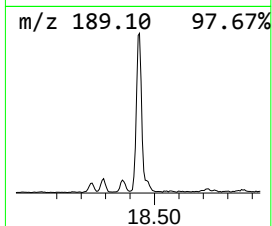
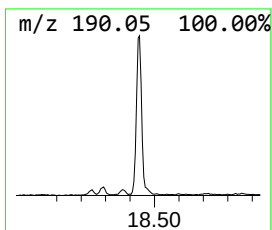
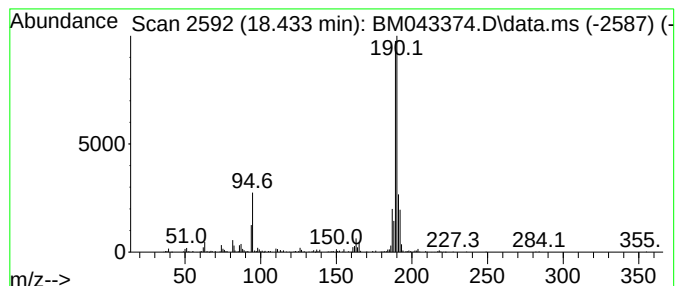
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 4H-Cyclopenta[def]phenanthrene Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.433	2.19 ng/ul	488803	Phenanthrene-d10	17.363

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	70
3			6H-Cyclobuta[jk]phenanthrene	190	C15H10	083469-43-6	59
4			4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	58
5			Methyl diselenide	190	C2H6Se2	007101-31-7	43



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM121523\
Data File : BM043374.D
Acq On : 16 Dec 2023 01:02
Operator : MA/JU
Sample : 05626-17
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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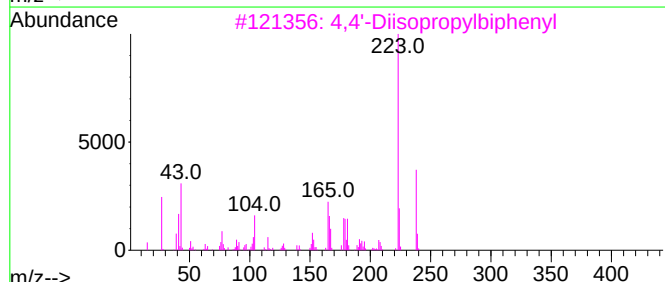
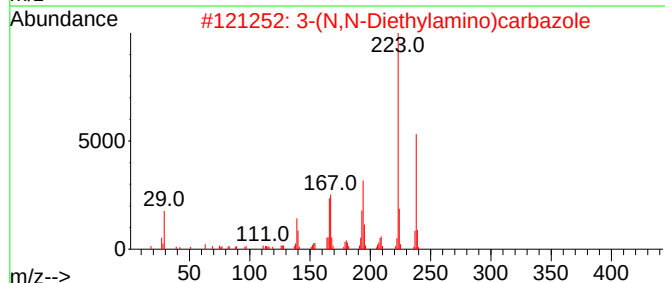
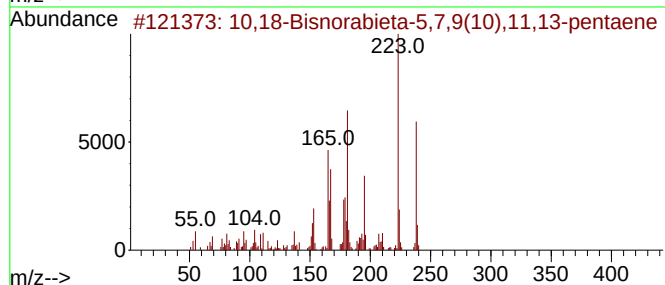
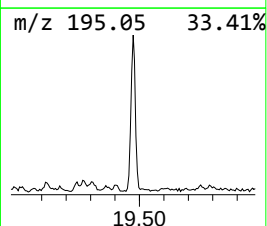
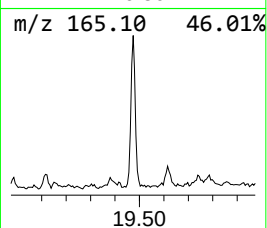
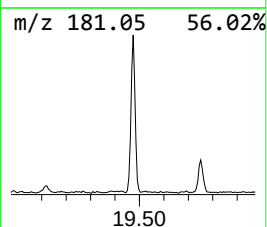
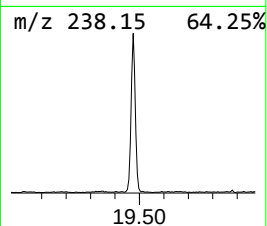
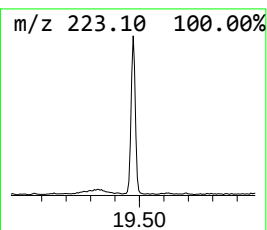
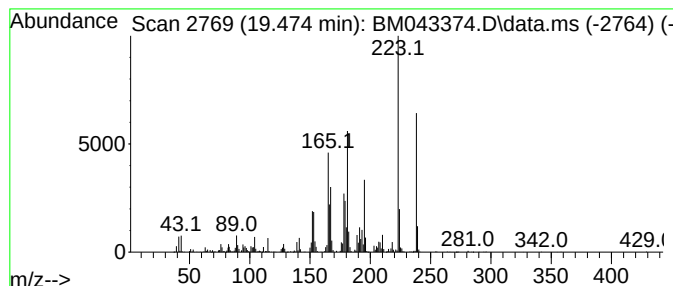
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 10,18-Bisnorabieta-5,7,9(10... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.474	3.06 ng/ul	881494	Chrysene-d12	21.539

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	10,18-Bisnorabieta-5,7,9(10),11,...	238	C18H22	006566-19-4	99		
2	3-(N,N-Diethylamino)carbazole	238	C16H18N2	054994-31-9	60		
3	4,4'-Diisopropylbiphenyl	238	C18H22	018970-30-4	55		
4	9-Acridinecarboxylic acid	223	C14H9NO2	005336-90-3	45		
5	4,4'-Diisopropylbiphenyl	238	C18H22	018970-30-4	45		



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM121523\
Data File : BM043374.D
Acq On : 16 Dec 2023 01:02
Operator : MA/JU
Sample : 05626-17
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
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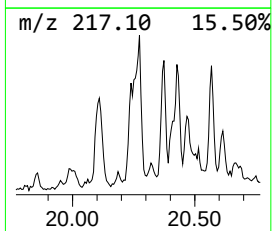
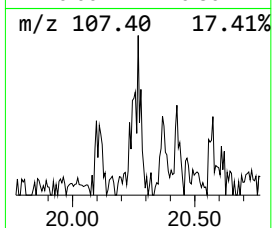
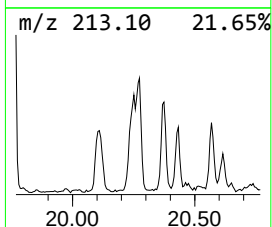
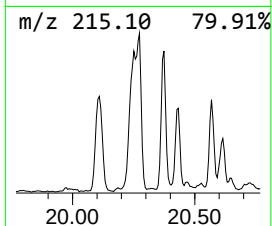
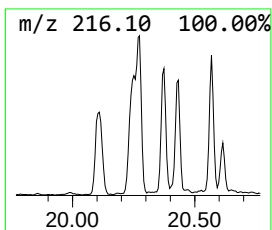
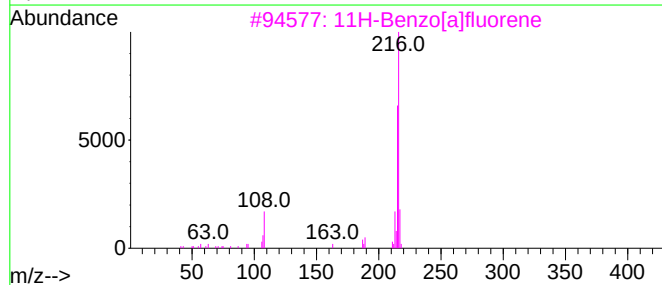
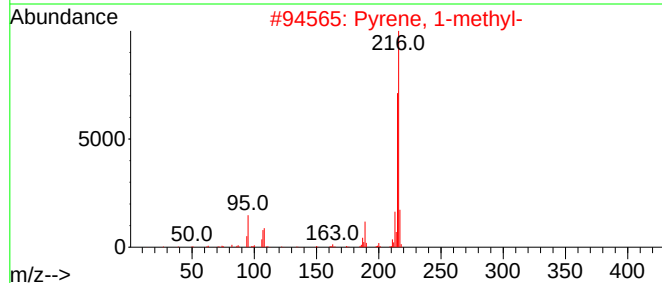
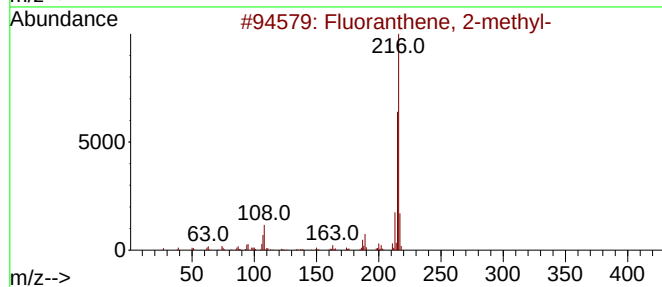
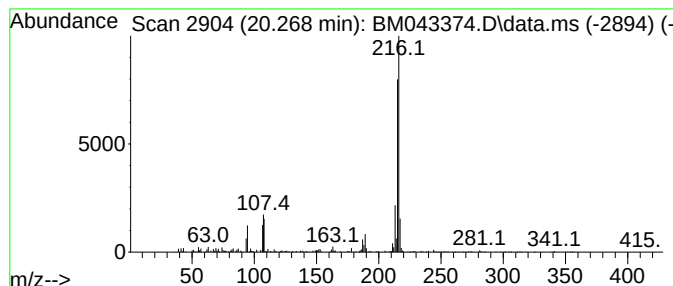
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 Fluoranthene, 2-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.268	3.15 ng/ul	907124	Chrysene-d12	21.539

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	94
2			Pyrene, 1-methyl-	216	C17H12	002381-21-7	93
3			11H-Benzo[a]fluorene	216	C17H12	000238-84-6	92
4			11H-Benzo[b]fluorene	216	C17H12	000243-17-4	91
5			11H-Benzo[b]fluorene	216	C17H12	000243-17-4	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM121523\
Data File : BM043374.D
Acq On : 16 Dec 2023 01:02
Operator : MA/JU
Sample : 05626-17
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Instrument :
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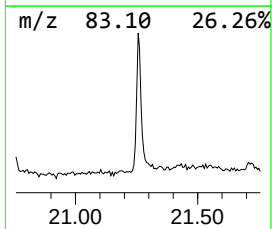
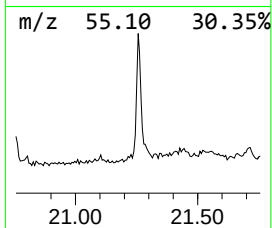
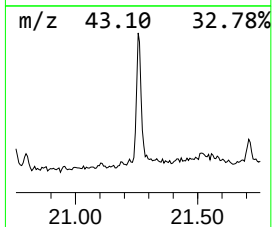
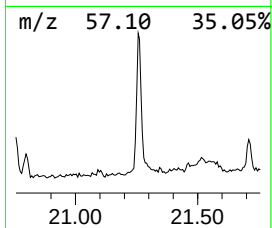
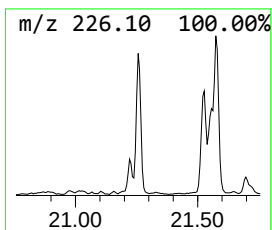
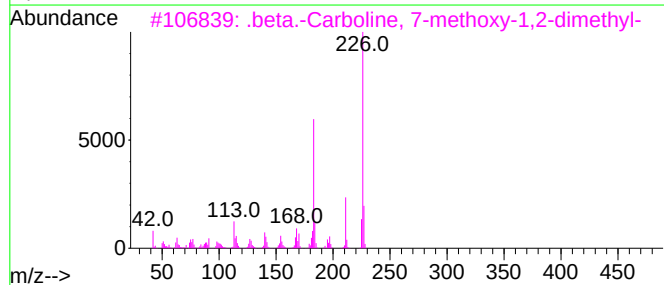
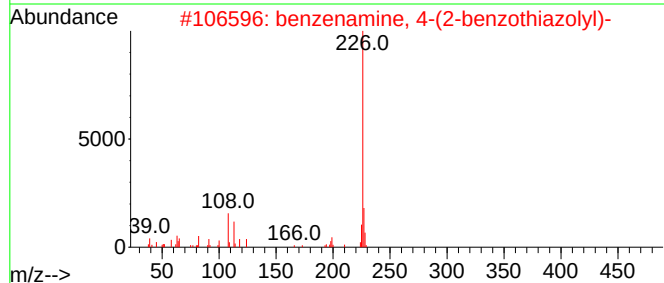
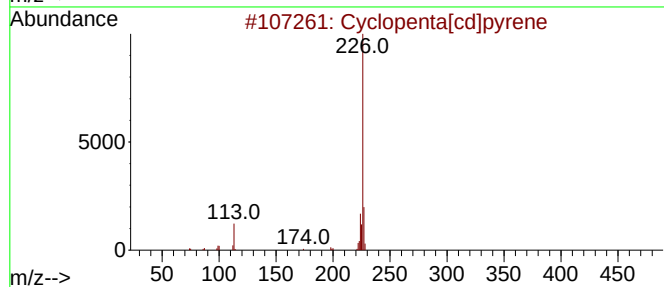
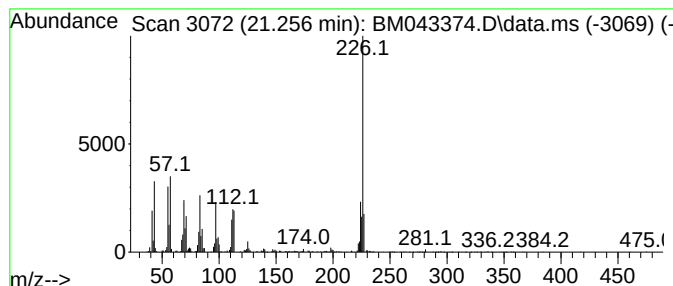
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 Cyclopenta[cd]pyrene Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.256	3.49 ng/ul	1006740	Chrysene-d12	21.539

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclopenta[cd]pyrene	226	C18H10	027208-37-3	60
2			benzenamine, 4-(2-benzothiazolyl)-	226	C13H10N2S	1000400-93-4	47
3			.beta.-Carboline, 7-methoxy-1,2-...	226	C14H14N2O	006519-18-2	43
4			Benzene, [4-(3-ethynylphenyl)-1,...	226	C18H10	1000115-87-3	38
5			Benzo[ghi]fluoranthene	226	C18H10	000203-12-3	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM121523\
Data File : BM043374.D
Acq On : 16 Dec 2023 01:02
Operator : MA/JU
Sample : 05626-17
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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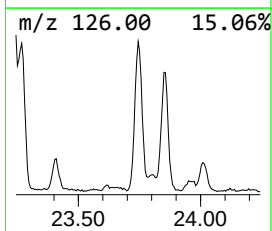
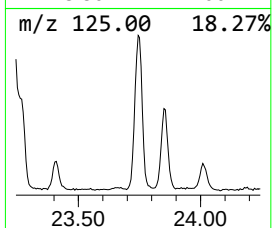
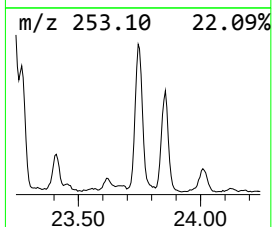
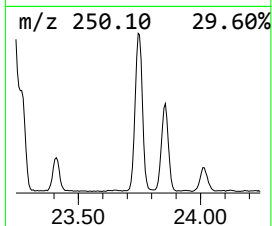
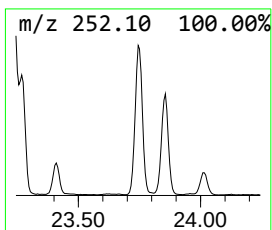
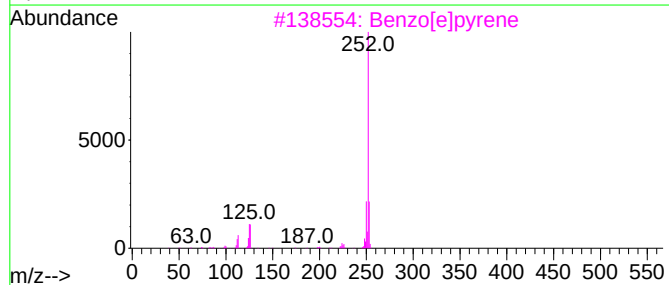
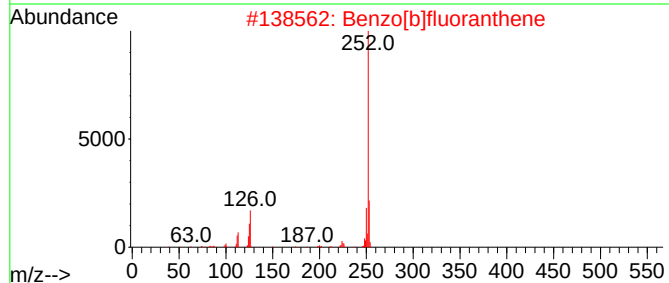
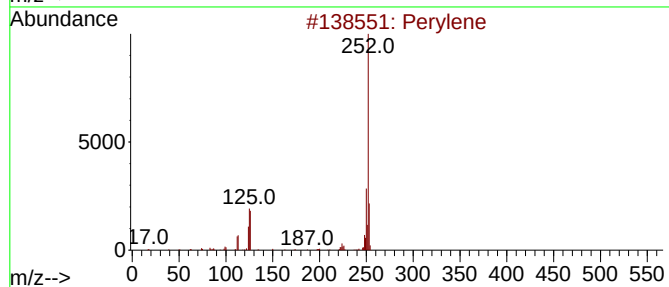
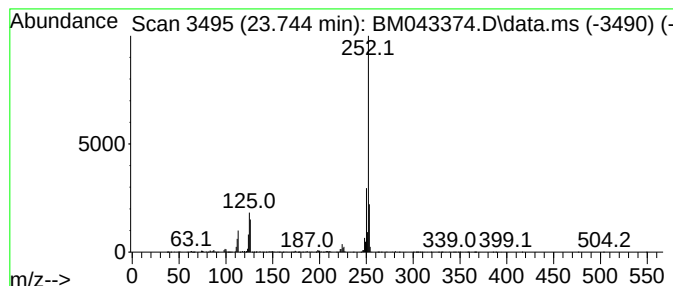
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 Perylene Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.744	11.14 ng/ul	1997080	Perylene-d12	23.962

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Perylene	252	C20H12	000198-55-0	99
2			Benzo[b]fluoranthene	252	C20H12	000205-99-2	98
3			Benzo[e]pyrene	252	C20H12	000192-97-2	98
4			Benzo[a]pyrene	252	C20H12	000050-32-8	96
5			Benzo[k]fluoranthene	252	C20H12	000207-08-9	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM121523\
Data File : BM043374.D
Acq On : 16 Dec 2023 01:02
Operator : MA/JU
Sample : 05626-17
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
DCLG3

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM121523.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST0.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
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2-Propanol, 1-(...	7.775	18.2	ng/ul	2528350	1	7.992	2778680	20.0
Aceto phenone	8.828	2.1	ng/ul	289784	1	7.992	2778680	20.0
n-Hexadecanoic ...	18.262	2.8	ng/ul	628481	4	17.363	4459770	20.0
4H-Cyclopenta[d...	18.433	2.2	ng/ul	488803	4	17.363	4459770	20.0
10,18-Bisnorabi...	19.474	3.1	ng/ul	881494	5	21.539	5761810	20.0
Fluoranthene, 2...	20.268	3.1	ng/ul	907124	5	21.539	5761810	20.0
Cyclopenta[cd]p...	21.256	3.5	ng/ul	1006740	5	21.539	5761810	20.0
Perylene	23.744	11.1	ng/ul	1997080	6	23.962	3584750	20.0