

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM013025\
 Data File : BM049529.D
 Acq On : 30 Jan 2025 16:23
 Operator : RC/JU
 Sample : Q1189-13DL2 30X
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 A44R4DL2

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-BM012825.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BM049529.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.546	431	435	444	rBV2	36099	63602	0.73%	0.087%
2	6.587	607	612	617	rBV4	11246	21563	0.25%	0.030%
3	6.675	622	627	632	rVV5	12854	22885	0.26%	0.031%
4	6.734	633	637	648	rVB4	21188	53126	0.61%	0.073%
5	6.869	653	660	666	rVV3	17509	31031	0.36%	0.043%
6	7.063	688	693	705	rVB3	25040	45703	0.53%	0.063%
7	7.181	705	713	721	rBV2	39126	84923	0.98%	0.117%
8	7.263	723	727	731	rVB4	14426	22352	0.26%	0.031%
9	7.510	762	769	778	rBV	1110344	1729241	19.87%	2.373%
10	7.769	806	813	828	rBV	94969	243028	2.79%	0.333%
11	8.040	854	859	867	rBV	65055	113845	1.31%	0.156%
12	8.245	890	894	901	rBV3	20389	38887	0.45%	0.053%
13	8.334	903	909	916	rVB	35549	65967	0.76%	0.091%
14	8.663	959	965	973	rVB3	16654	33619	0.39%	0.046%
15	8.775	979	984	990	rBV3	15577	28997	0.33%	0.040%
16	9.263	1062	1067	1072	rBV3	10674	17113	0.20%	0.023%
17	9.375	1079	1086	1093	rVB6	15924	32154	0.37%	0.044%
18	9.716	1136	1144	1150	rBV2	30784	61646	0.71%	0.085%
19	9.781	1150	1155	1159	rBV2	18818	32087	0.37%	0.044%
20	9.857	1165	1168	1176	rVB2	19796	37406	0.43%	0.051%
21	9.934	1176	1181	1189	rVB3	20166	42024	0.48%	0.058%
22	10.275	1230	1239	1256	rVB	1378886	2342532	26.91%	3.214%
23	11.322	1410	1417	1419	rBV7	11673	20332	0.23%	0.028%
24	11.504	1444	1448	1454	rVB6	12888	28897	0.33%	0.040%
25	11.610	1461	1466	1471	rBV7	9980	17088	0.20%	0.023%
26	11.804	1486	1499	1506	rBV3	64357	193692	2.23%	0.266%
27	11.951	1517	1524	1533	rVB5	13886	32505	0.37%	0.045%
28	12.169	1556	1561	1567	rBV4	13439	27613	0.32%	0.038%
29	12.692	1644	1650	1655	rBV8	13193	26957	0.31%	0.037%
30	12.922	1683	1689	1692	rBV7	13027	20776	0.24%	0.029%
31	12.992	1696	1701	1708	rVB10	7127	18189	0.21%	0.025%
32	13.127	1718	1724	1728	rBV3	26275	55620	0.64%	0.076%
33	13.392	1764	1769	1776	rBV	41932	75264	0.86%	0.103%
34	13.475	1779	1783	1788	rVV2	32951	55122	0.63%	0.076%
35	13.522	1788	1791	1795	rVV3	14867	21130	0.24%	0.029%
36	13.575	1795	1800	1802	rVV2	68033	101036	1.16%	0.139%

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Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM012825.MA.M
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37	13.604	1802	1805	1811	rVV	131308	215956	2.48%	0.296%
38	13.657	1811	1814	1818	rVV2	19180	30031	0.35%	0.041%
39	13.710	1818	1823	1826	rVB3	25203	41411	0.48%	0.057%
40	13.874	1841	1851	1865	rBV	476444	872076	10.02%	1.197%
41	13.980	1865	1869	1874	rVB5	19806	26469	0.30%	0.036%
42	14.157	1892	1899	1905	rBV	2046203	2961754	34.03%	4.064%
43	14.216	1906	1909	1917	rVB	65822	104869	1.20%	0.144%
44	14.374	1930	1936	1940	rBV8	22026	48378	0.56%	0.066%
45	14.474	1946	1953	1962	rBV9	30145	102595	1.18%	0.141%
46	14.563	1966	1968	1971	rVB3	19861	22881	0.26%	0.031%
47	14.639	1978	1981	1985	rVB	29942	37962	0.44%	0.052%
48	14.698	1987	1991	1998	rBV9	21039	34420	0.40%	0.047%
49	14.774	1998	2004	2009	rBV3	62836	157414	1.81%	0.216%
50	14.869	2017	2020	2025	rVB2	21656	30620	0.35%	0.042%
51	14.933	2027	2031	2037	rBV5	27103	61157	0.70%	0.084%
52	15.033	2043	2048	2055	rVB	112526	166156	1.91%	0.228%
53	15.121	2059	2063	2065	rBV2	30166	42388	0.49%	0.058%
54	15.157	2065	2069	2074	rVB	135313	187152	2.15%	0.257%
55	15.216	2074	2079	2080	rBV5	65380	98200	1.13%	0.135%
56	15.339	2096	2100	2104	rVB2	58252	72957	0.84%	0.100%
57	15.421	2109	2114	2115	rBV2	46377	75334	0.87%	0.103%
58	15.627	2146	2149	2151	rBV2	20168	20399	0.23%	0.028%
59	15.892	2189	2194	2197	rBV3	81920	144862	1.66%	0.199%
60	15.927	2197	2200	2204	rVB	75902	105645	1.21%	0.145%
61	16.033	2214	2218	2220	rBV	39399	50193	0.58%	0.069%
62	16.216	2244	2249	2255	rBV4	51567	128989	1.48%	0.177%
63	16.274	2256	2259	2262	rVV2	34137	44680	0.51%	0.061%
64	16.492	2292	2296	2300	rVB2	80556	117398	1.35%	0.161%
65	16.568	2306	2309	2316	rVB3	55427	89843	1.03%	0.123%
66	16.645	2319	2322	2328	rBV4	51622	93356	1.07%	0.128%
67	16.910	2361	2367	2371	rBV	2420695	3341661	38.39%	4.585%
68	16.951	2371	2374	2379	rVV	477621	611548	7.03%	0.839%
69	17.010	2380	2384	2385	rVV	88459	111689	1.28%	0.153%
70	17.039	2385	2389	2395	rVB	630231	871344	10.01%	1.196%
71	17.104	2395	2400	2407	rBV3	119944	217211	2.50%	0.298%
72	17.321	2434	2437	2440	rBV	70875	89550	1.03%	0.123%
73	17.433	2454	2456	2461	rVB2	83117	93612	1.08%	0.128%
74	17.786	2512	2516	2520	rBV	138881	196087	2.25%	0.269%
75	17.833	2520	2524	2531	rVB	286002	392496	4.51%	0.539%
76	17.915	2533	2538	2543	rBV2	243084	348601	4.00%	0.478%

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77	17.980	2544	2549	2553	rBV2	648239	1088142	12.50%	1.493%
78	18.633	2657	2660	2665	rVB2	147798	201098	2.31%	0.276%
79	18.715	2670	2674	2680	rBV	295895	445531	5.12%	0.611%
80	18.780	2682	2685	2689	rVB2	200607	266153	3.06%	0.365%
81	18.857	2694	2698	2706	rBV	649052	1184811	13.61%	1.626%
82	18.980	2714	2719	2725	rVB	5079094	6427580	73.84%	8.819%
83	19.133	2741	2745	2749	rBV	296647	392401	4.51%	0.538%
84	19.345	2772	2781	2790	rBV	5255622	8704482	100.00%	11.943%
85	19.427	2791	2795	2800	rBV3	225585	497879	5.72%	0.683%
86	19.527	2810	2812	2818	rVB	295943	390468	4.49%	0.536%
87	19.680	2834	2838	2849	rVB3	545054	1212393	13.93%	1.664%
88	19.951	2881	2884	2888	rVB	381349	439932	5.05%	0.604%
89	20.009	2890	2894	2898	rVB2	861997	1191764	13.69%	1.635%
90	20.151	2914	2918	2921	rBV	679620	873706	10.04%	1.199%
91	20.192	2922	2925	2930	rVB	677676	923496	10.61%	1.267%
92	20.603	2992	2995	3001	rBV	552967	727111	8.35%	0.998%
93	20.809	3027	3030	3033	rVB	496590	553767	6.36%	0.760%
94	20.845	3033	3036	3044	rBV2	516405	1129991	12.98%	1.550%
95	21.139	3080	3086	3091	rBV2	3806117	8558646	98.32%	11.743%
96	21.192	3091	3095	3103	rVB	2569687	3958915	45.48%	5.432%
97	21.797	3195	3198	3202	rVB	859003	1104295	12.69%	1.515%
98	23.068	3408	3414	3420	rBV2	2722081	7058034	81.09%	9.684%
99	23.803	3534	3539	3550	rVB	1427308	3160786	36.31%	4.337%
100	23.939	3555	3562	3571	rBV	1569275	4072751	46.79%	5.588%

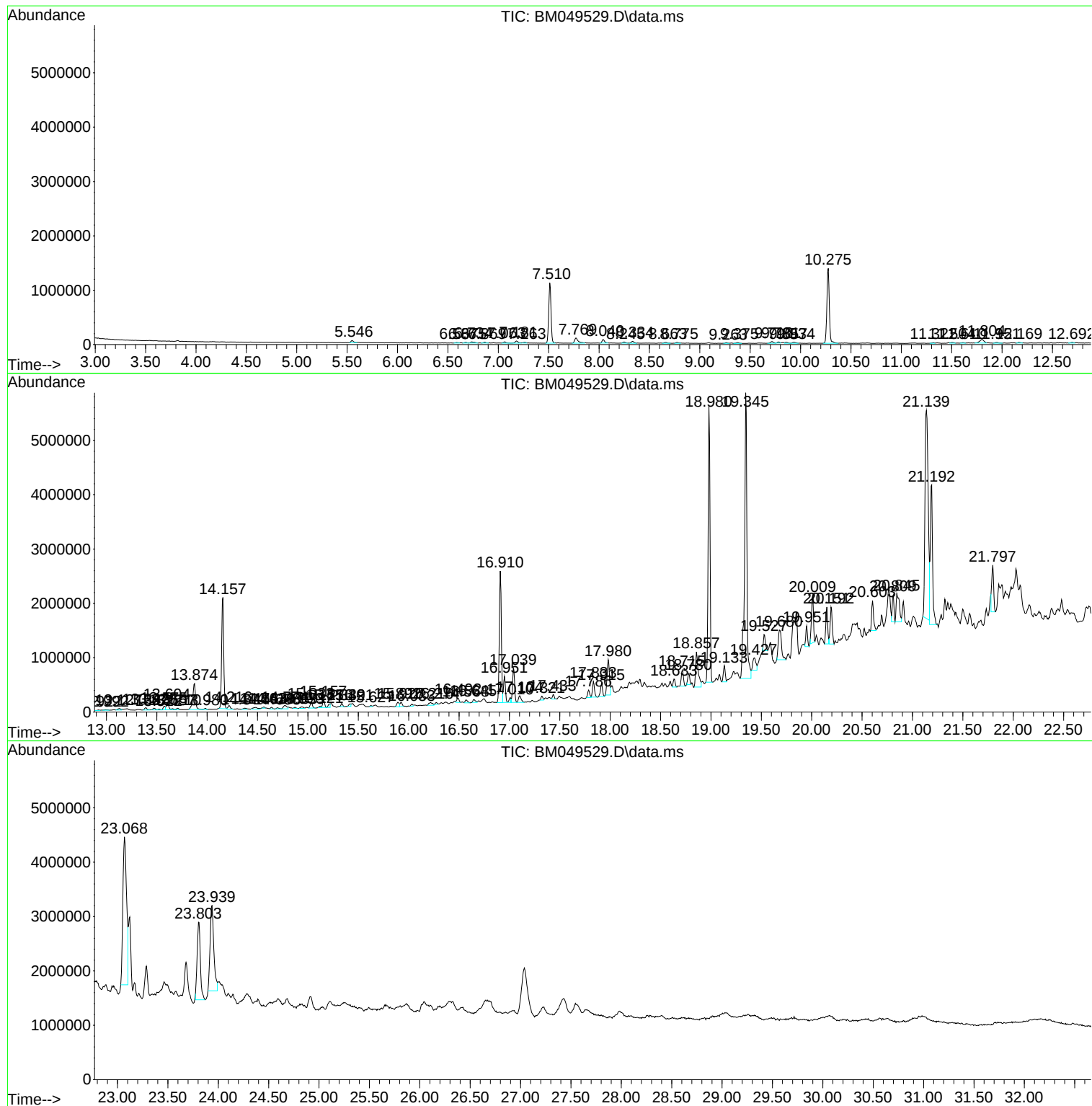
Sum of corrected areas: 72881428

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM012825.MA.M
Quant Title : SVOA CALIBRATION

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



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Misc :
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Instrument :
BNA_M
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A44R4DL2

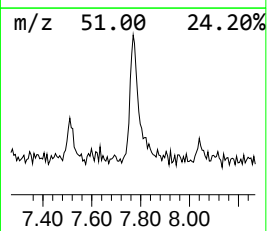
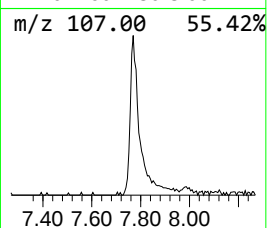
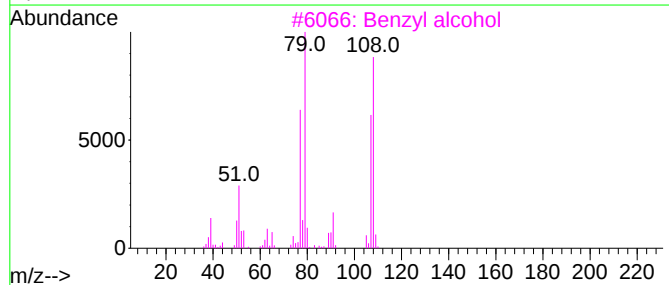
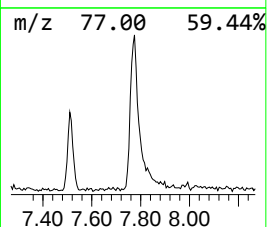
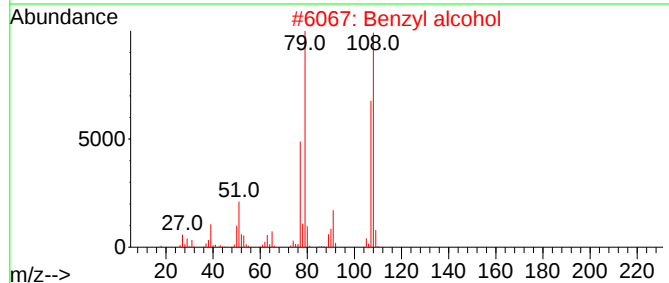
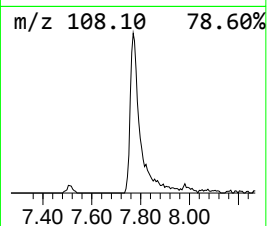
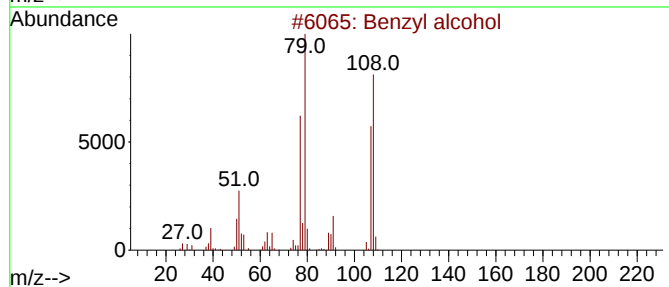
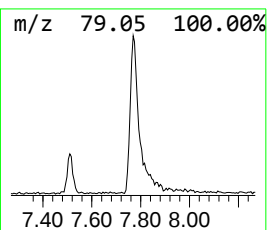
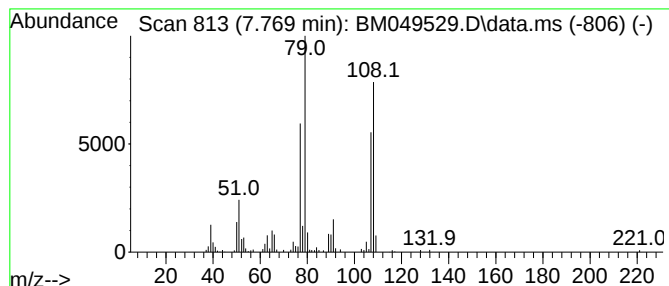
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM012825.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Benzyl alcohol Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.769	2.81 ng/ul	243028	1,4-Dichlorobenzene-d4	7.510

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzyl alcohol	108	C7H8O	000100-51-6	95
2			Benzyl alcohol	108	C7H8O	000100-51-6	93
3			Benzyl alcohol	108	C7H8O	000100-51-6	93
4			N-Cbz-glycylglycine p-nitropheny...	387	C18H17N3O7	013574-81-7	91
5			N-.alpha.,N-.omega.-Di-cbz-L-arg...	442	C22H26N4O6	053934-75-1	91



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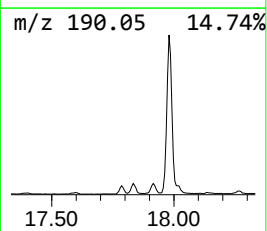
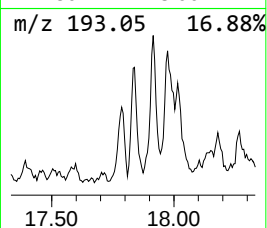
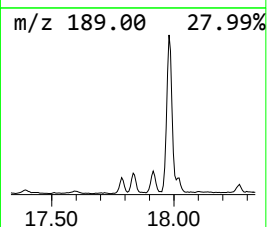
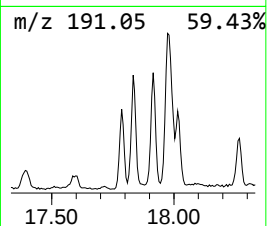
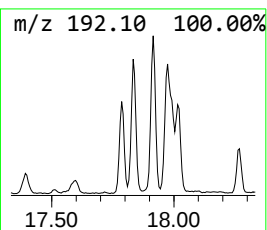
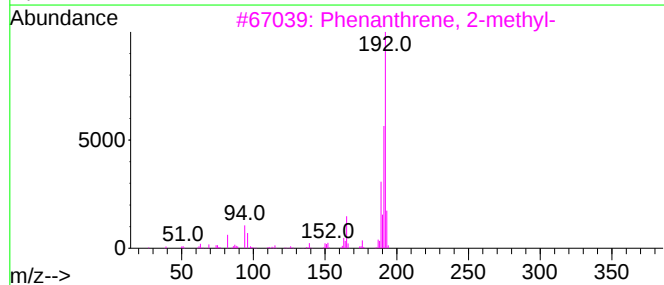
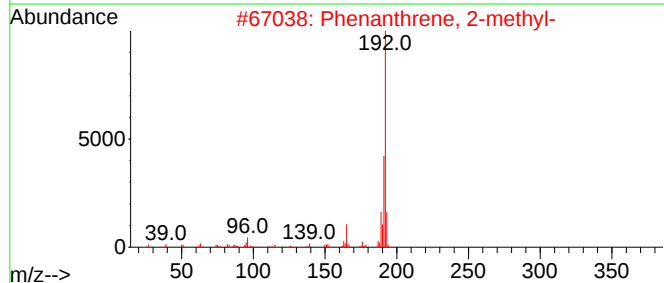
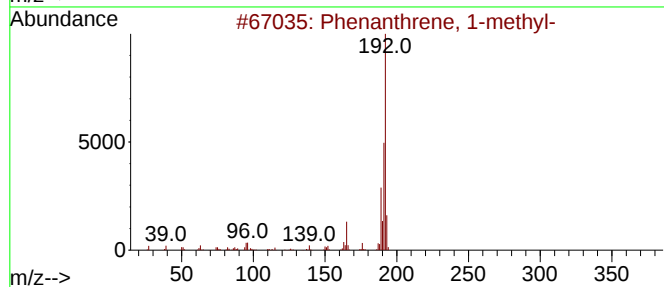
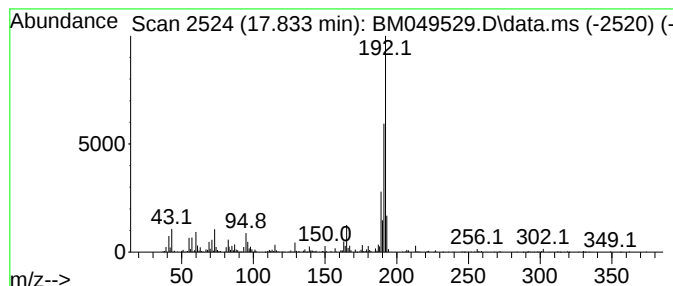
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM012825.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 Phenanthrene, 1-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.833	2.35 ng/ul	392496	Phenanthrene-d10	16.910

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	96
2			Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	95
3			Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	94
4			Phenanthrene, 4-methyl-	192	C15H12	000832-64-4	92
5			Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	91



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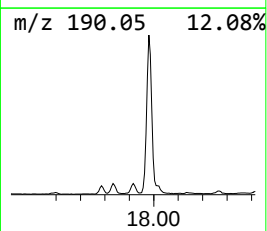
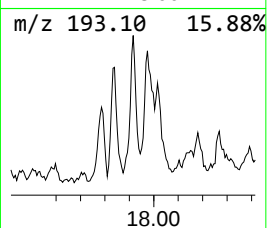
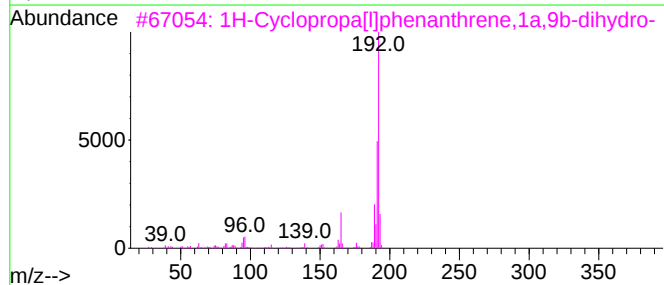
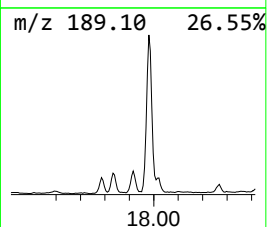
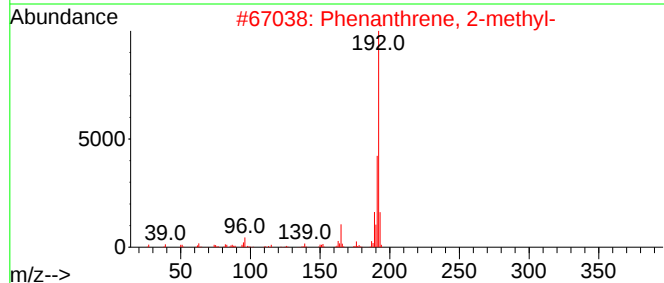
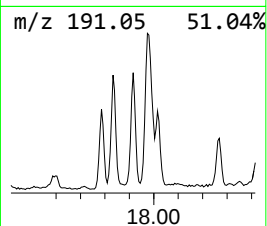
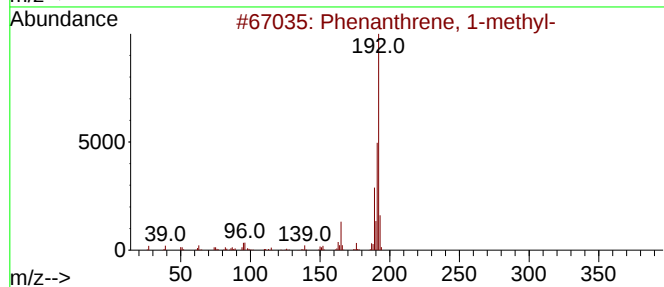
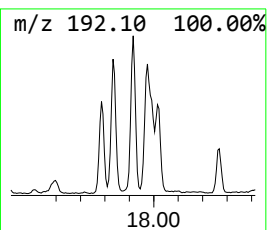
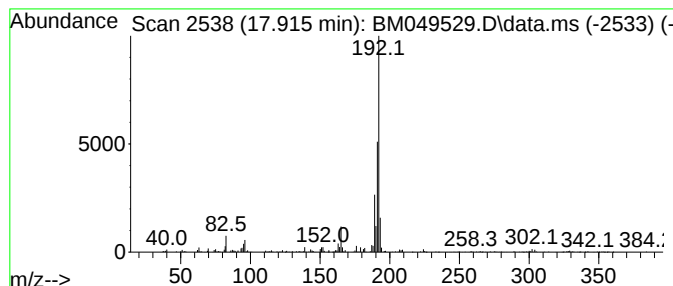
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Quant Title : SVOA CALIBRATION

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

Peak Number 3 Phenanthrene, 2-methyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.915	2.09 ng/ul	348601	Phenanthrene-d10	16.910

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	97
2			Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	97
3			1H-Cyclopropa[1]phenanthrene,1a,...	192	C15H12	000949-41-7	97
4			Anthracene, 9-methyl-	192	C15H12	000779-02-2	95
5			Anthracene, 9-methyl-	192	C15H12	000779-02-2	95



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM013025\
Data File : BM049529.D
Acq On : 30 Jan 2025 16:23
Operator : RC/JU
Sample : Q1189-13DL2 30X
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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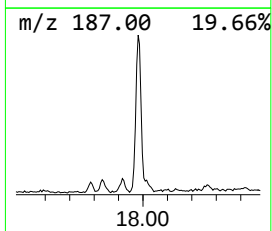
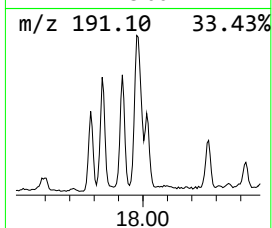
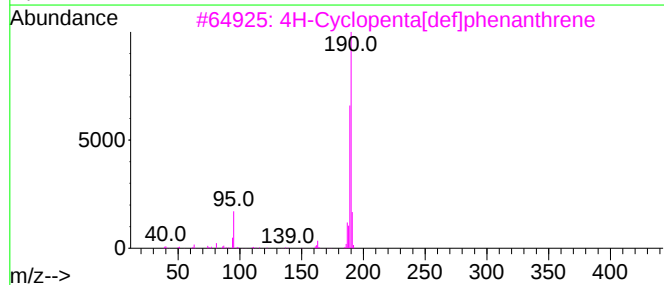
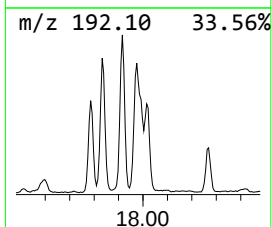
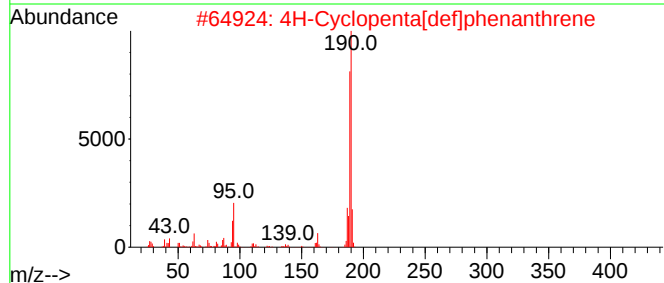
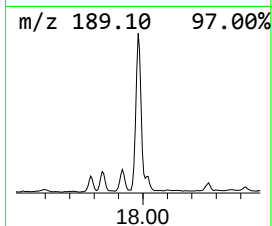
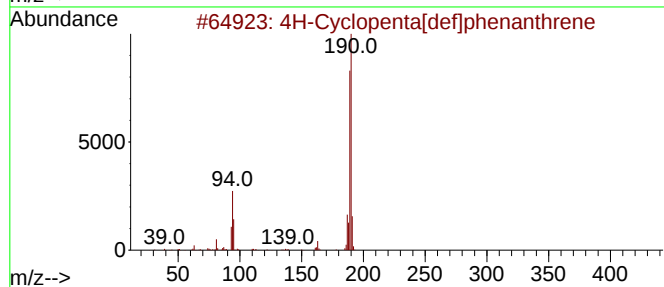
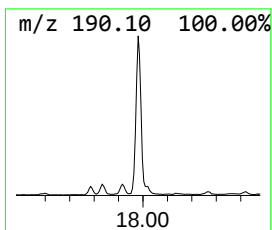
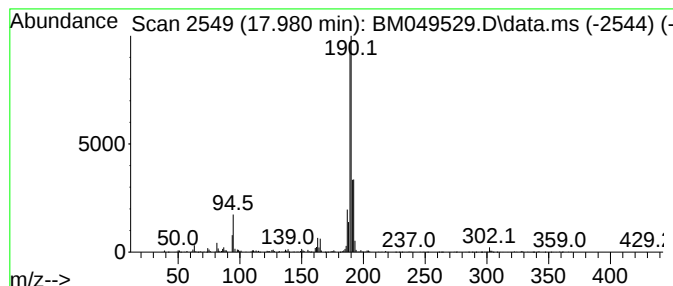
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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.980	6.51 ng/ul	1088140	Phenanthrene-d10	16.910

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	94
2			4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	76
3			4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	68
4			2,3,5,6-Tetramethylterephthalald...	190	C12H14O2	007072-01-7	50
5			6H-Cyclobuta[jk]phenanthrene	190	C15H10	083469-43-6	50



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM013025\
Data File : BM049529.D
Acq On : 30 Jan 2025 16:23
Operator : RC/JU
Sample : Q1189-13DL2 30X
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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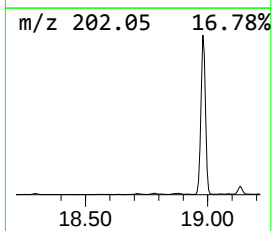
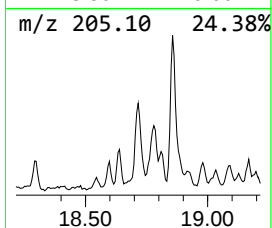
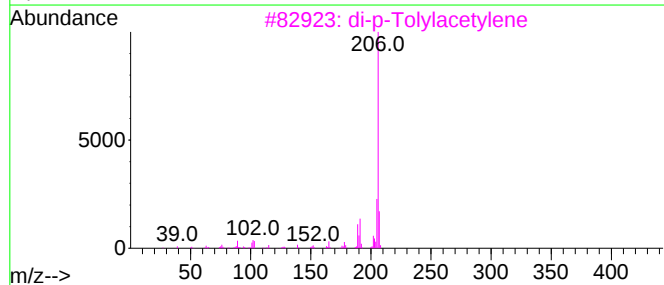
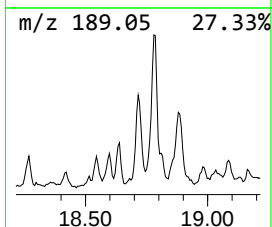
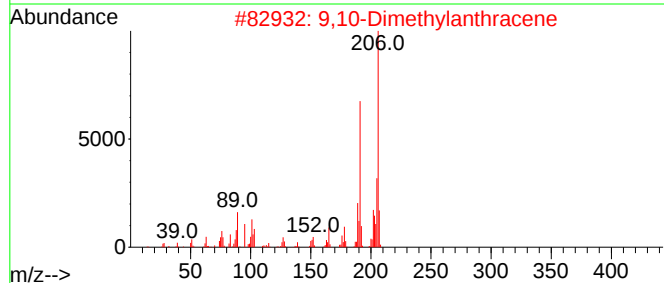
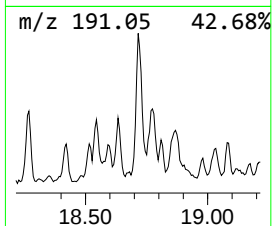
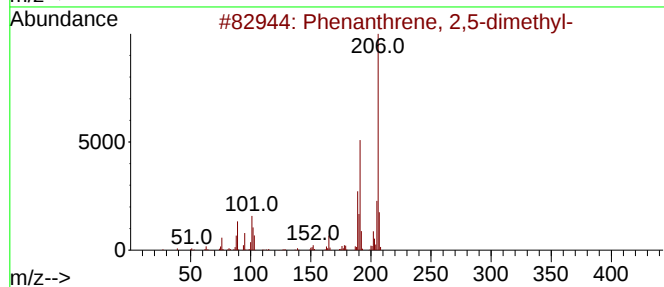
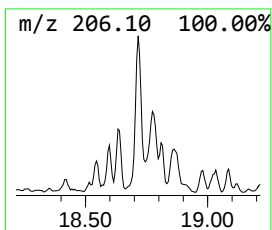
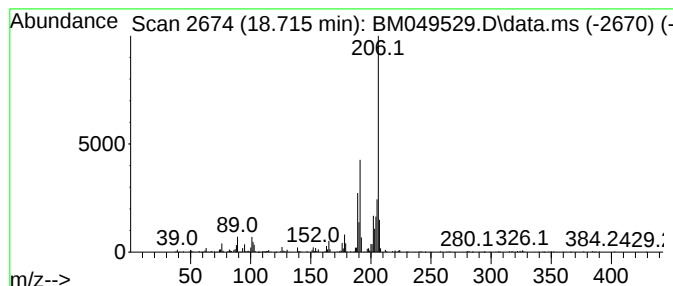
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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.715	2.67 ng/ul	445531	Phenanthrene-d10	16.910

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenanthrene, 2,5-dimethyl-	206	C16H14	003674-66-6	96
2			9,10-Dimethylantracene	206	C16H14	000781-43-1	95
3			di-p-Tolylacetylene	206	C16H14	002789-88-0	94
4			9,10-Dimethylantracene	206	C16H14	000781-43-1	91
5			Phenanthrene, 3,6-dimethyl-	206	C16H14	001576-67-6	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM013025\
Data File : BM049529.D
Acq On : 30 Jan 2025 16:23
Operator : RC/JU
Sample : Q1189-13DL2 30X
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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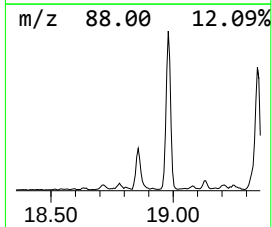
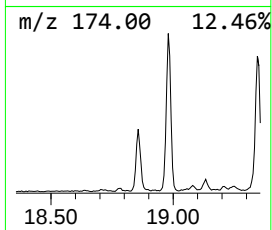
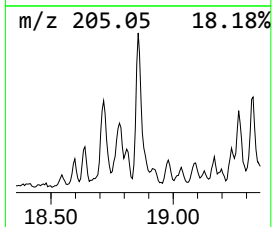
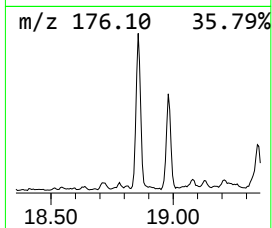
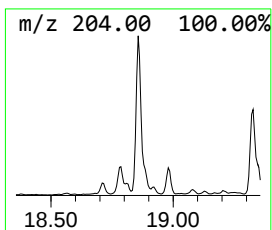
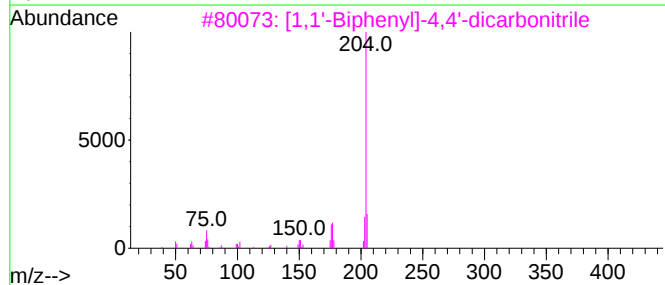
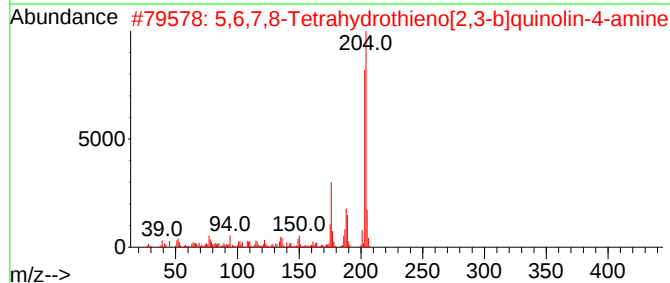
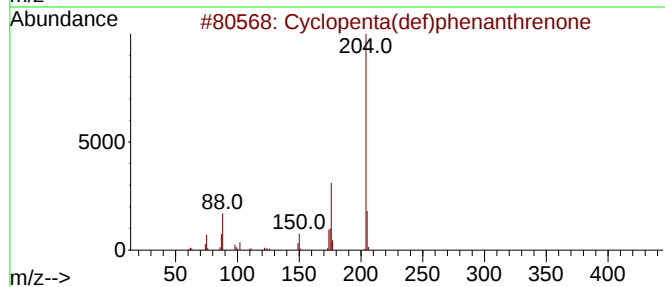
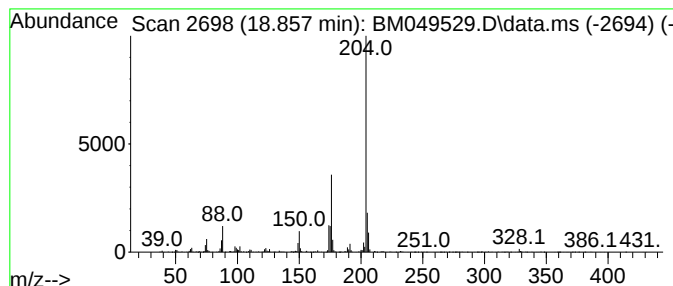
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 Cyclopenta(def)phenanthrenone Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.857	7.09 ng/ul	1184810	Phenanthrene-d10	16.910

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclopenta(def)phenanthrenone	204	C15H8O	005737-13-3	96
2			5,6,7,8-Tetrahydrothieno[2,3-b]q...	204	C11H12N2S	122914-50-5	72
3			[1,1'-Biphenyl]-4,4'-dicarbonitrile	204	C14H8N2	001591-30-6	59
4			(E)-Ethyl 3-oxo-2-(pyrid-2-yl)me...	219	C12H13NO3	1000147-59-7	59
5			[1,1'-Biphenyl]-2,2'-dicarbonitrile	204	C14H8N2	004341-02-0	59



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM013025\
Data File : BM049529.D
Acq On : 30 Jan 2025 16:23
Operator : RC/JU
Sample : Q1189-13DL2 30X
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
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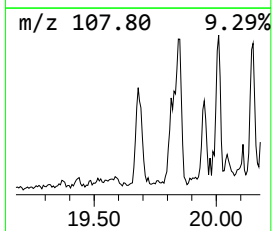
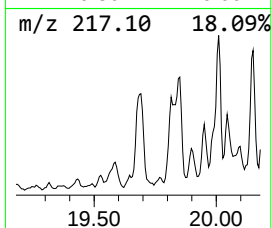
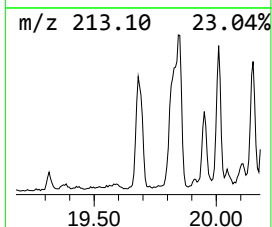
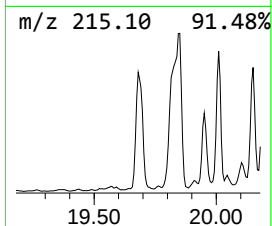
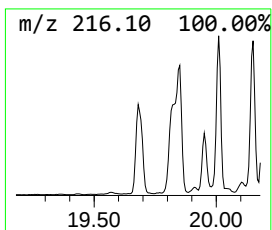
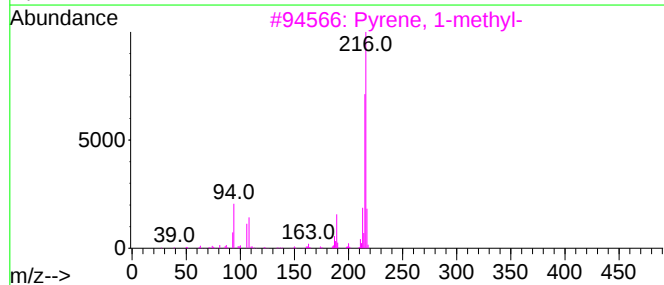
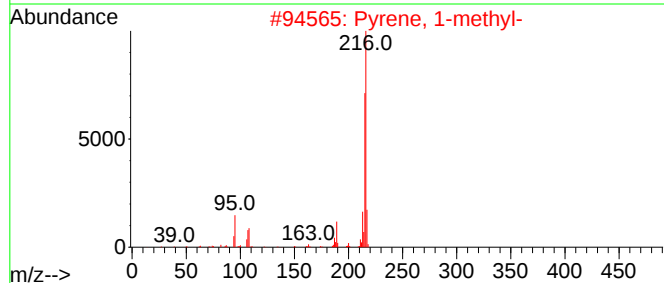
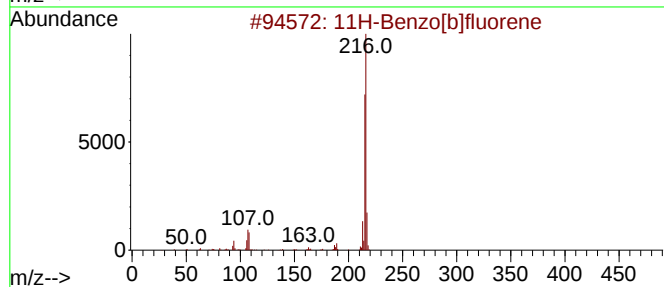
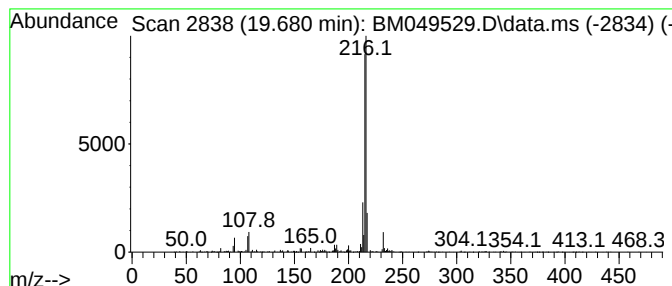
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 11H-Benzo[b]fluorene Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.680	2.83 ng/ul	1212390	Chrysene-d12	21.151

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			11H-Benzo[b]fluorene	216	C17H12	000243-17-4	87
2			Pyrene, 1-methyl-	216	C17H12	002381-21-7	83
3			Pyrene, 1-methyl-	216	C17H12	002381-21-7	74
4			Pyrene, 1-methyl-	216	C17H12	002381-21-7	64
5			11H-Benzo[a]fluorene	216	C17H12	000238-84-6	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM013025\
Data File : BM049529.D
Acq On : 30 Jan 2025 16:23
Operator : RC/JU
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Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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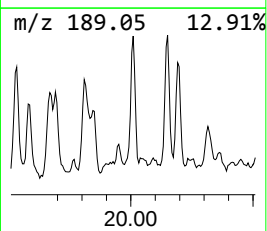
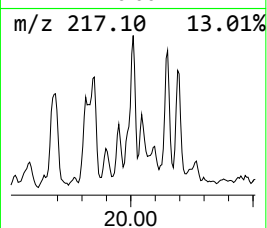
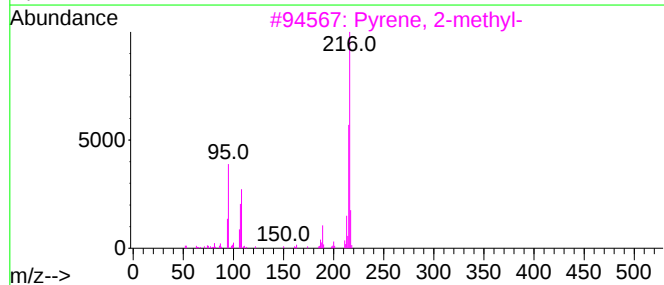
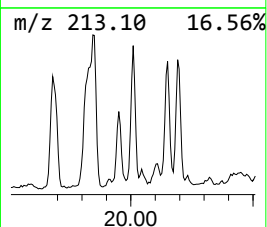
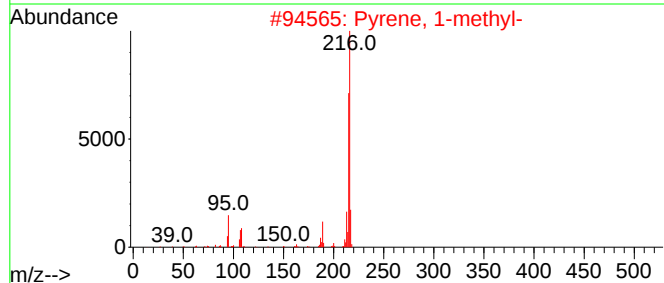
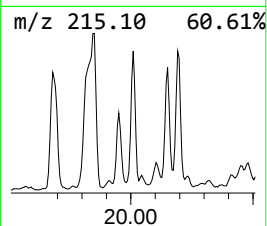
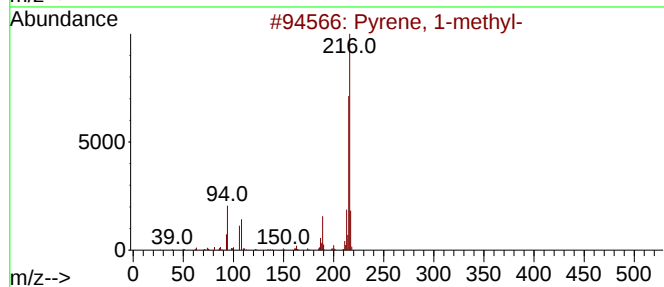
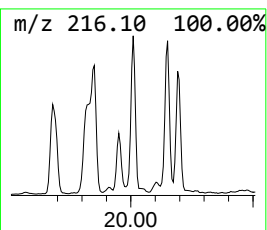
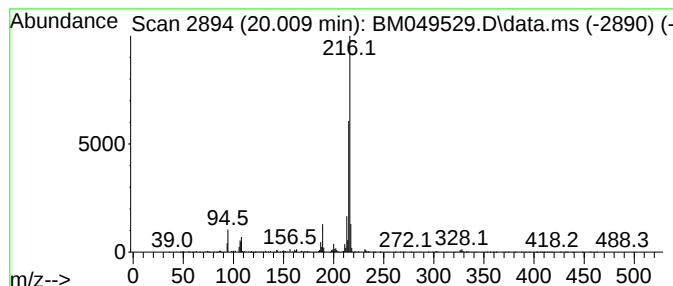
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 Pyrene, 1-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.009	2.78 ng/ul	1191760	Chrysene-d12	21.151

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pyrene, 1-methyl-	216	C17H12	002381-21-7	96
2			Pyrene, 1-methyl-	216	C17H12	002381-21-7	94
3			Pyrene, 2-methyl-	216	C17H12	003442-78-2	94
4			Pyrene, 4-methyl-	216	C17H12	003353-12-6	93
5			Fluoranthene, 2-methyl-	216	C17H12	003543-31-6	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM013025\
Data File : BM049529.D
Acq On : 30 Jan 2025 16:23
Operator : RC/JU
Sample : Q1189-13DL2 30X
Misc :
ALS Vial : 10 Sample Multiplier: 1

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ClientSampleId :
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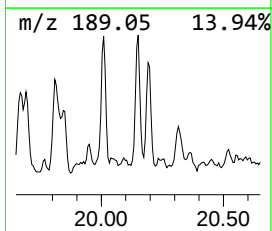
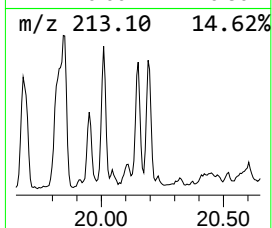
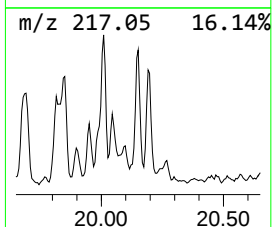
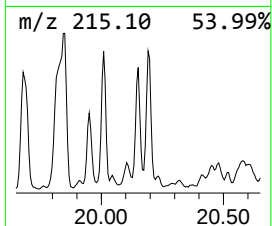
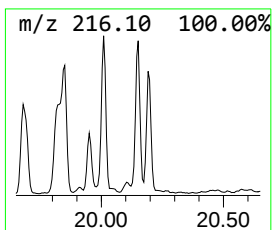
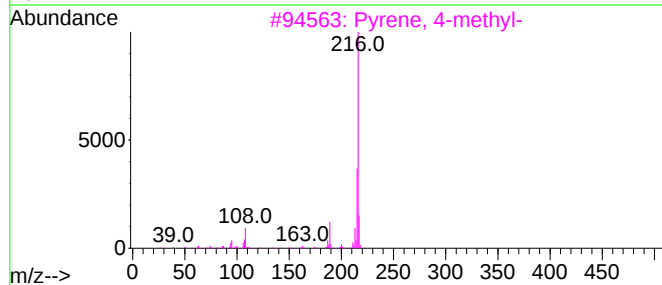
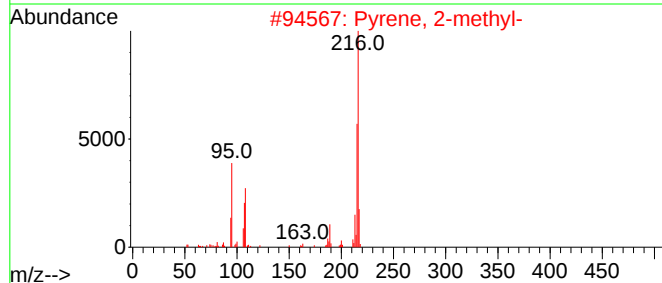
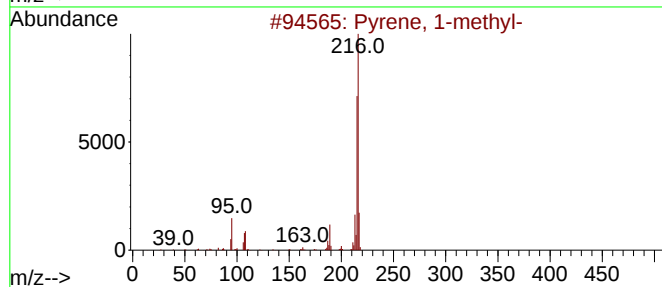
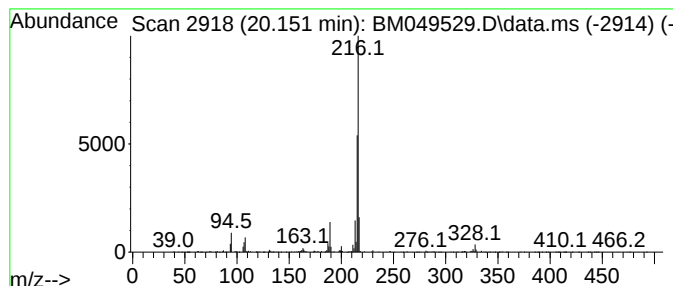
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 Pyrene, 2-methyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.151	2.04 ng/ul	873706	Chrysene-d12	21.151

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pyrene, 1-methyl-	216	C17H12	002381-21-7	96
2			Pyrene, 2-methyl-	216	C17H12	003442-78-2	94
3			Pyrene, 4-methyl-	216	C17H12	003353-12-6	94
4			Pyrene, 1-methyl-	216	C17H12	002381-21-7	93
5			Pyrene, 1-methyl-	216	C17H12	002381-21-7	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM013025\
Data File : BM049529.D
Acq On : 30 Jan 2025 16:23
Operator : RC/JU
Sample : Q1189-13DL2 30X
Misc :
ALS Vial : 10 Sample Multiplier: 1

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ClientSampleId :
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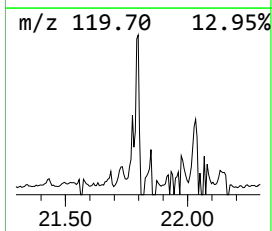
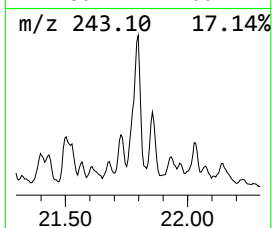
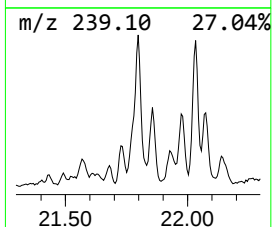
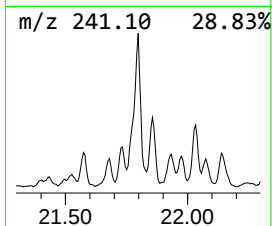
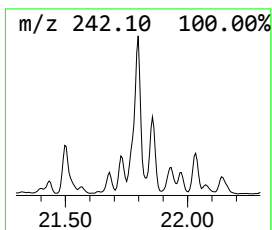
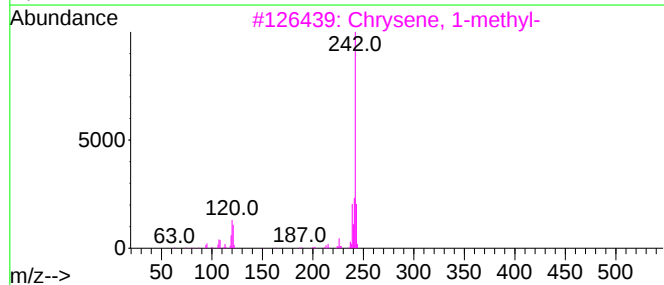
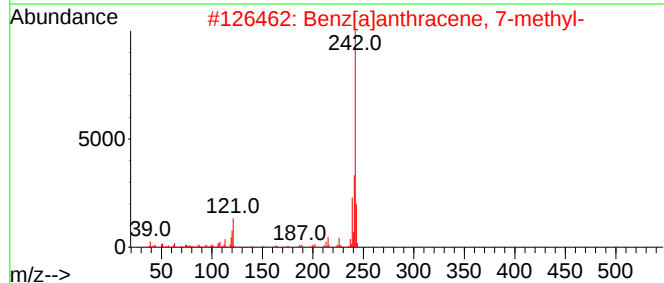
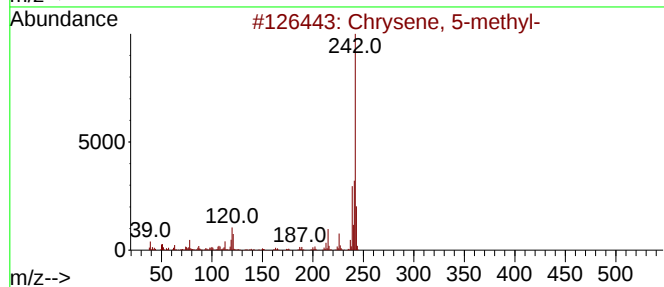
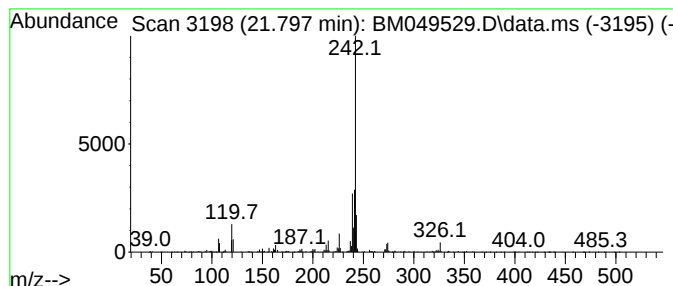
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM012825.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 12 Chrysene, 5-methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.797	2.58 ng/ul	1104300	Chrysene-d12	21.151

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Chrysene, 5-methyl-	242	C19H14	003697-24-3	93
2		Benz[a]anthracene, 7-methyl-	242	C19H14	002541-69-7	93
3		Chrysene, 1-methyl-	242	C19H14	003351-28-8	90
4		2-Methylchrysene	242	C19H14	003351-32-4	89
5		Chrysene, 1-methyl-	242	C19H14	003351-28-8	89



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM013025\
Data File : BM049529.D
Acq On : 30 Jan 2025 16:23
Operator : RC/JU
Sample : Q1189-13DL2 30X
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
A44R4DL2

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM012825.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
Benzyl alcohol	7.769	2.8	ng/u1	243028	1	7.510	1729240	20.0
Phenanthrene, 1...	17.833	2.4	ng/u1	392496	4	16.910	3341660	20.0
Phenanthrene, 2...	17.915	2.1	ng/u1	348601	4	16.910	3341660	20.0
4H-Cyclopenta[d...	17.980	6.5	ng/u1	1088140	4	16.910	3341660	20.0
Phenanthrene, 2...	18.715	2.7	ng/u1	445531	4	16.910	3341660	20.0
Cyclopenta(def)...	18.857	7.1	ng/u1	1184810	4	16.910	3341660	20.0
11H-Benzo[b]flu...	19.680	2.8	ng/u1	1212390	5	21.151	8558650	20.0
Pyrene, 1-methyl-	20.009	2.8	ng/u1	1191760	5	21.151	8558650	20.0
Pyrene, 2-methyl-	20.151	2.0	ng/u1	873706	5	21.151	8558650	20.0
Chrysene, 5-met...	21.797	2.6	ng/u1	1104300	5	21.151	8558650	20.0