

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM010224\
 Data File : BM043702.D
 Acq On : 02 Jan 2024 19:54
 Operator : MA/JU
 Sample : 05920-07
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 GCF58

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: BM043702.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.369	27	31	38	rVB	68695	94499	1.29%	0.087%
2	3.793	100	103	118	rBV	721952	1197521	16.36%	1.102%
3	3.904	119	122	129	rVB2	37375	57574	0.79%	0.053%
4	5.310	357	361	368	rVV	28799	50862	0.69%	0.047%
5	5.522	390	397	404	rBV5	24326	54040	0.74%	0.050%
6	5.992	470	477	483	rBV4	28644	59252	0.81%	0.055%
7	6.622	573	584	591	rVB	223353	356415	4.87%	0.328%
8	7.128	662	670	682	rBV	1310077	2096710	28.64%	1.929%
9	7.245	684	690	694	rVV	63412	103823	1.42%	0.096%
10	7.298	694	699	708	rVV	884194	1363187	18.62%	1.254%
11	7.381	708	713	723	rVB	50065	100472	1.37%	0.092%
12	7.487	723	731	738	rBV	1190920	1878681	25.66%	1.728%
13	7.957	802	811	818	rBV	843133	1360486	18.58%	1.251%
14	8.169	840	847	853	rBV	326085	498679	6.81%	0.459%
15	8.675	926	933	943	rVV	1557504	2426390	33.15%	2.232%
16	9.133	1004	1011	1025	rBV	1124772	1876817	25.64%	1.726%
17	9.710	1098	1109	1117	rBV5	26234	74711	1.02%	0.069%
18	9.851	1126	1133	1140	rBV	967531	1617746	22.10%	1.488%
19	10.075	1164	1171	1180	rVV6	54631	114604	1.57%	0.105%
20	10.392	1217	1225	1241	rBV	1448581	2556111	34.92%	2.351%
21	10.539	1244	1250	1260	rVB2	45680	91034	1.24%	0.084%
22	10.763	1281	1288	1297	rVB	1065802	1862656	25.44%	1.713%
23	10.916	1307	1314	1335	rVV	560234	1146240	15.66%	1.054%
24	11.322	1375	1383	1392	rBV3	25579	68208	0.93%	0.063%
25	12.121	1510	1519	1523	rBV4	34365	77875	1.06%	0.072%
26	12.339	1548	1556	1564	rBV3	92095	172937	2.36%	0.159%
27	13.098	1675	1685	1694	rBV10	20917	52271	0.71%	0.048%
28	13.368	1727	1731	1735	rVV7	23578	46970	0.64%	0.043%
29	13.468	1743	1748	1750	rVV2	67652	108934	1.49%	0.100%
30	13.492	1750	1752	1756	rVV2	54508	74631	1.02%	0.069%
31	13.598	1765	1770	1775	rVV3	66183	102583	1.40%	0.094%
32	13.721	1786	1791	1795	rBV7	27833	50265	0.69%	0.046%
33	13.768	1795	1799	1806	rBV5	48438	91597	1.25%	0.084%
34	13.857	1808	1814	1818	rBV	707857	1069269	14.61%	0.984%
35	13.904	1818	1822	1832	rVB2	271842	433844	5.93%	0.399%
36	14.015	1832	1841	1849	rBV	2436973	3402705	46.48%	3.130%

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Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM121523.MA.M
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37	14.110	1852	1857	1863	rVB2	190994	293975	4.02%	0.270%
38	14.292	1881	1888	1896	rVV	2878211	4344879	59.35%	3.997%
39	14.368	1897	1901	1904	rVV5	26887	51381	0.70%	0.047%
40	14.410	1904	1908	1913	rVV4	49551	96276	1.32%	0.089%
41	14.480	1913	1920	1925	rVB4	103558	211949	2.90%	0.195%
42	14.545	1925	1931	1934	rBV	112577	163983	2.24%	0.151%
43	14.598	1934	1940	1946	rVV	1507008	2422272	33.09%	2.228%
44	14.810	1969	1976	1991	rBV2	1227890	2485508	33.95%	2.286%
45	14.968	1998	2003	2007	rBV2	115970	176375	2.41%	0.162%
46	15.010	2007	2010	2014	rVV5	40461	62109	0.85%	0.057%
47	15.074	2018	2021	2026	rVB	69097	94465	1.29%	0.087%
48	15.262	2048	2053	2060	rVB5	29067	65039	0.89%	0.060%
49	15.586	2102	2108	2117	rBV	3649506	5250652	71.73%	4.830%
50	15.709	2124	2129	2135	rBV	1016482	1421041	19.41%	1.307%
51	15.821	2144	2148	2152	rBV4	69238	98404	1.34%	0.091%
52	15.874	2152	2157	2163	rVB	610751	826501	11.29%	0.760%
53	16.004	2172	2179	2183	rBV8	33776	74500	1.02%	0.069%
54	16.233	2212	2218	2223	rBV9	27025	53946	0.74%	0.050%
55	16.304	2224	2230	2235	rBV9	29785	66976	0.91%	0.062%
56	16.374	2238	2242	2248	rVB4	34543	53662	0.73%	0.049%
57	16.439	2248	2253	2257	rBV4	62759	105795	1.45%	0.097%
58	16.745	2299	2305	2310	rBV	101829	177873	2.43%	0.164%
59	16.992	2343	2347	2356	rVB6	41135	73182	1.00%	0.067%
60	17.239	2384	2389	2396	rBV3	53621	80892	1.11%	0.074%
61	17.339	2401	2406	2411	rVV	1841061	2662678	36.37%	2.449%
62	17.380	2411	2413	2417	rVV	128923	154317	2.11%	0.142%
63	17.439	2417	2423	2433	rVB2	3546190	5087628	69.50%	4.680%
64	17.551	2438	2442	2447	rVV6	36872	67518	0.92%	0.062%
65	17.703	2464	2468	2469	rVV4	51489	55695	0.76%	0.051%
66	17.727	2469	2472	2483	rVB3	93084	155254	2.12%	0.143%
67	18.121	2534	2539	2545	rBV3	52585	95496	1.30%	0.088%
68	18.186	2545	2550	2554	rBV	163466	271354	3.71%	0.250%
69	18.245	2555	2560	2574	rVV	1521535	2358554	32.22%	2.169%
70	18.356	2576	2579	2586	rVB	110158	180290	2.46%	0.166%
71	18.539	2606	2610	2612	rBV4	46950	73508	1.00%	0.068%
72	18.574	2612	2616	2623	rVB2	115751	201376	2.75%	0.185%
73	18.668	2629	2632	2638	rVV	75250	91049	1.24%	0.084%
74	18.727	2640	2642	2650	rVB3	57530	78293	1.07%	0.072%
75	19.198	2719	2722	2727	rVB	197364	240246	3.28%	0.221%
76	19.297	2735	2739	2744	rBV3	47149	67749	0.93%	0.062%

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77	19.368	2745	2751	2754	rBV3	160147	319240	4.36%	0.294%
78	19.397	2754	2756	2758	rBV	68982	74367	1.02%	0.068%
79	19.521	2773	2777	2787	rBV2	478882	817450	11.17%	0.752%
80	19.633	2792	2796	2800	rVB	198995	238432	3.26%	0.219%
81	19.733	2807	2813	2820	rBV	4712159	6268632	85.63%	5.766%
82	19.903	2838	2842	2848	rBV4	113241	173516	2.37%	0.160%
83	19.986	2854	2856	2860	rVB2	85588	87061	1.19%	0.080%
84	20.250	2897	2901	2904	rBV2	145900	198062	2.71%	0.182%
85	20.439	2928	2933	2942	rBV2	650038	1058841	14.46%	0.974%
86	20.609	2957	2962	2966	rBV	3233593	3845473	52.53%	3.537%
87	20.650	2966	2969	2976	rVB2	1570864	2246060	30.68%	2.066%
88	20.921	3010	3015	3038	rVB	4582905	7320403	100.00%	6.734%
89	21.239	3065	3069	3074	rVB	1528437	1761428	24.06%	1.620%
90	21.333	3081	3085	3096	rVB4	404957	839993	11.47%	0.773%
91	21.521	3113	3117	3130	rBV	1998724	2883588	39.39%	2.652%
92	22.103	3213	3216	3220	rBV	304926	376319	5.14%	0.346%
93	22.174	3222	3228	3240	rVB	1838435	3001239	41.00%	2.761%
94	23.227	3403	3407	3410	rBV	703234	1029723	14.07%	0.947%
95	23.274	3410	3415	3424	rVB	1981262	3496591	47.77%	3.216%
96	23.780	3495	3501	3510	rVB2	2864754	5918332	80.85%	5.444%
97	23.938	3522	3528	3537	rVB	1494126	3016273	41.20%	2.774%
98	24.603	3636	3641	3649	rVB	1084825	2282508	31.18%	2.100%
99	24.697	3653	3657	3668	rVB2	503659	1075668	14.69%	0.989%
100	27.403	4111	4117	4131	rVB3	984735	3002528	41.02%	2.762%

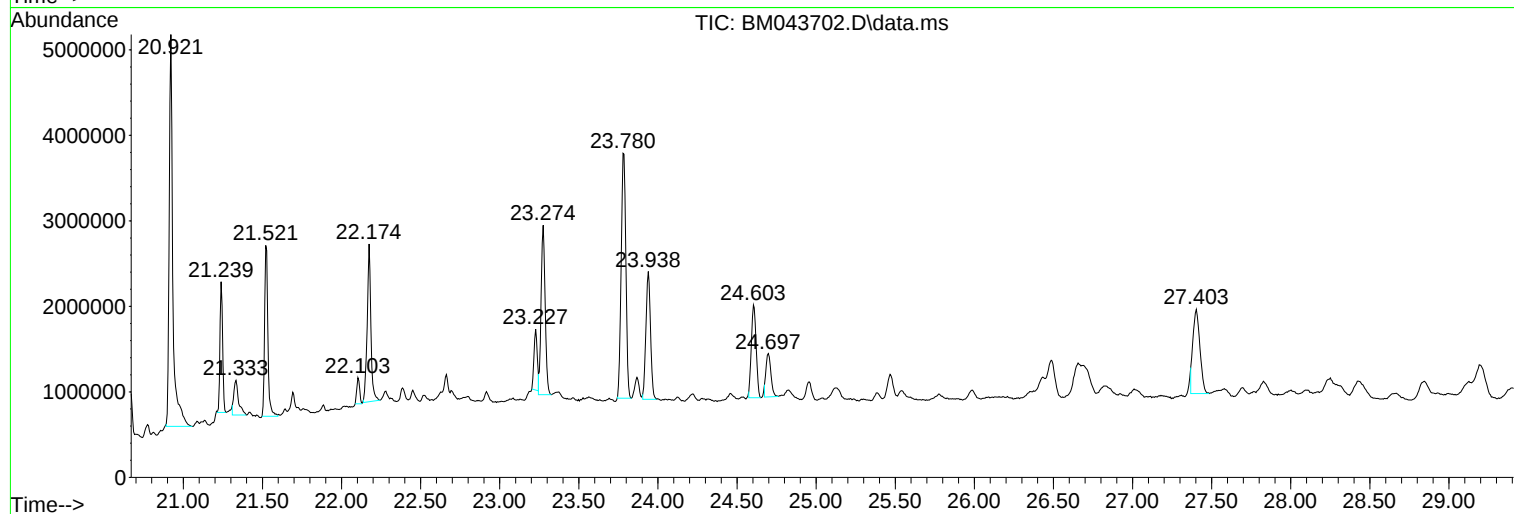
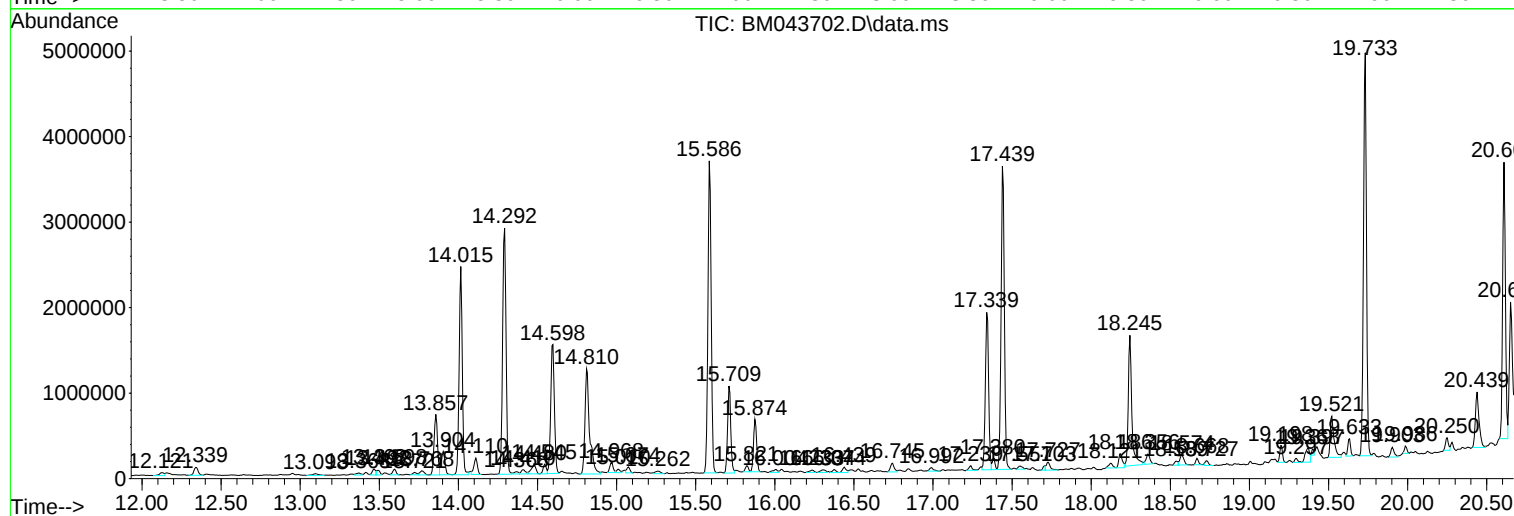
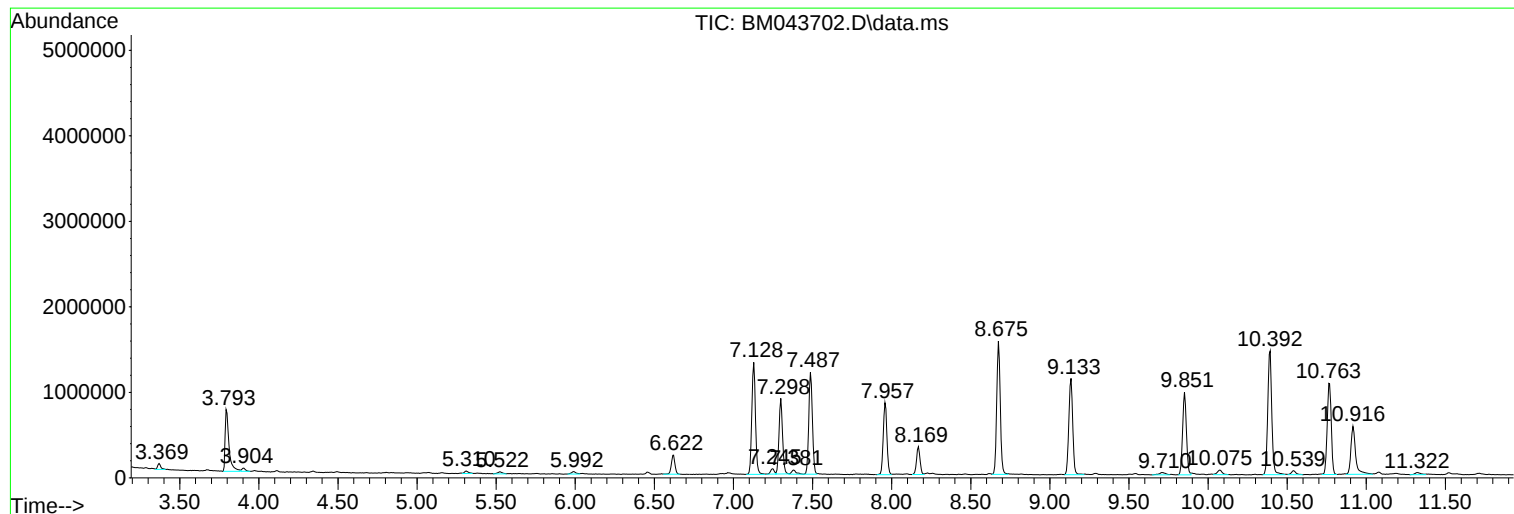
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM121523.MA.M
Quant Title : SVOA CALIBRATION

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



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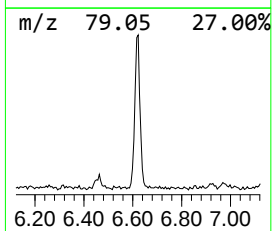
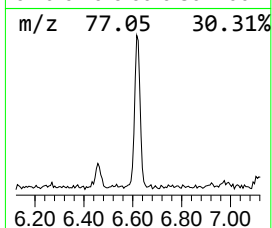
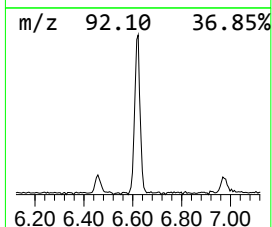
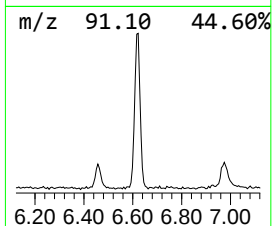
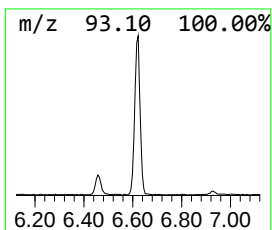
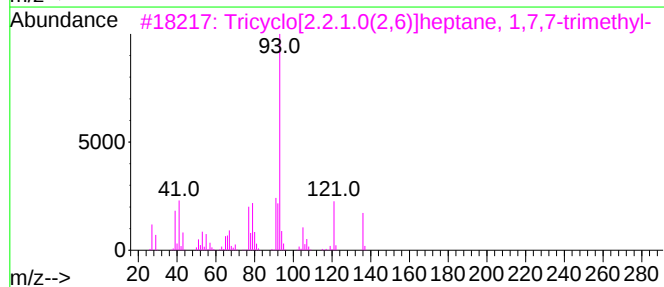
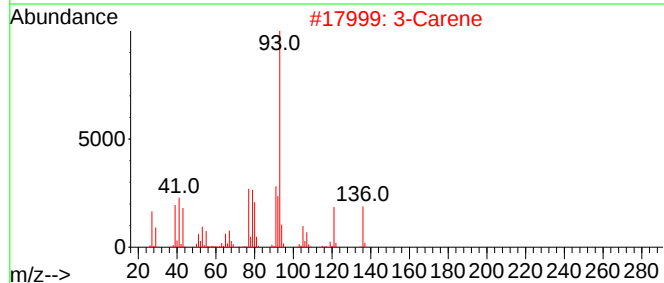
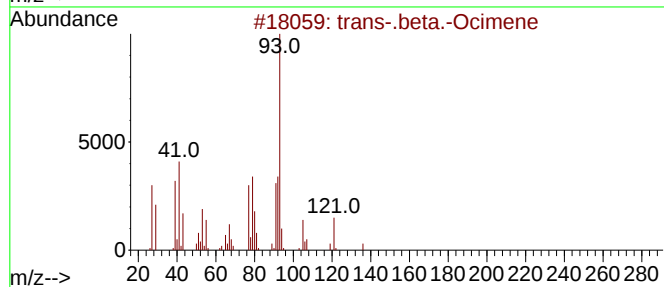
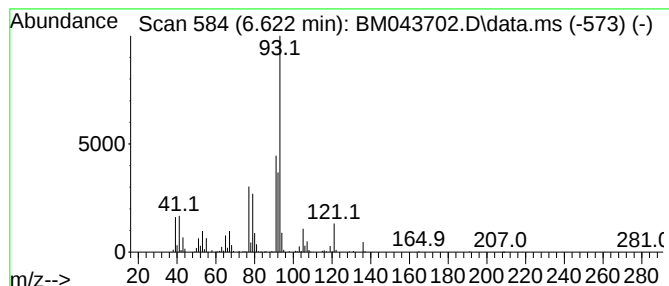
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM121523.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 trans-.beta.-Ocimene Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.622	5.24 ng/ul	356415	1,4-Dichlorobenzene-d4	7.957

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			trans-.beta.-Ocimene	136	C10H16	003779-61-1	94
2			3-Carene	136	C10H16	013466-78-9	91
3			Tricyclo[2.2.1.0(2,6)]heptane, 1...	136	C10H16	000508-32-7	91
4			(+)-3-Carene	136	C10H16	000498-15-7	91
5			Bicyclo[3.1.1]hept-2-ene, 3,6,6-...	136	C10H16	004889-83-2	91



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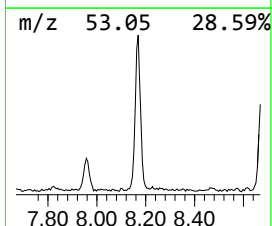
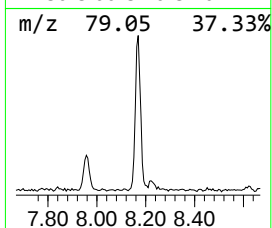
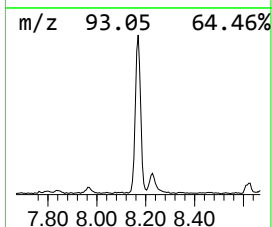
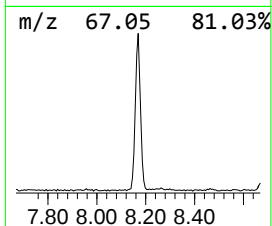
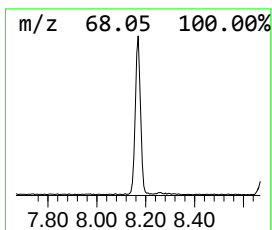
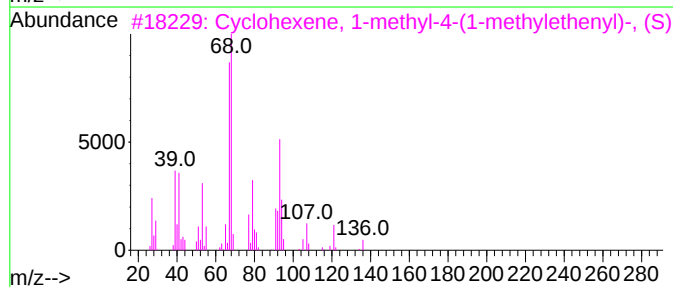
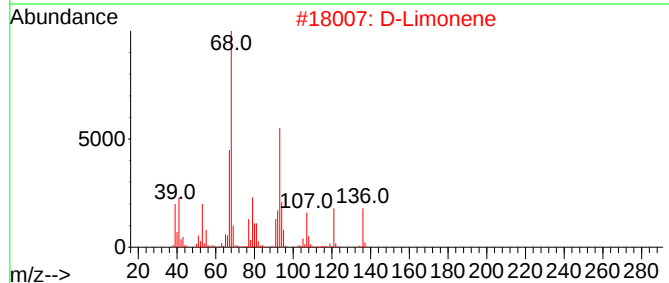
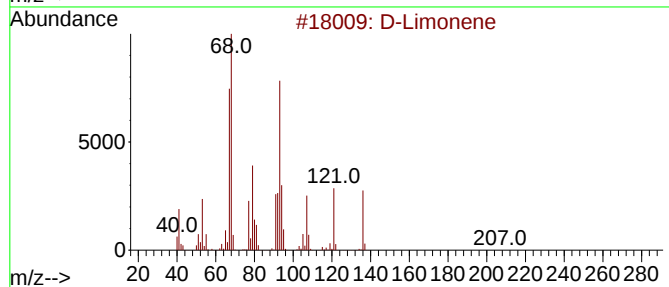
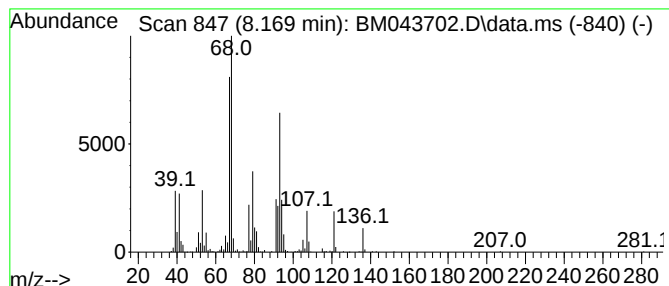
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM121523.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 D-Limonene Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.169	7.33 ng/ul	498679	1,4-Dichlorobenzene-d4	7.957

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	D-Limonene			136	C10H16	005989-27-5	98
2	D-Limonene			136	C10H16	005989-27-5	91
3	Cyclohexene, 1-methyl-4-(1-methy...			136	C10H16	005989-54-8	90
4	Limonene			136	C10H16	000138-86-3	87
5	D-Limonene			136	C10H16	005989-27-5	87



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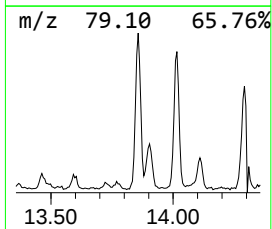
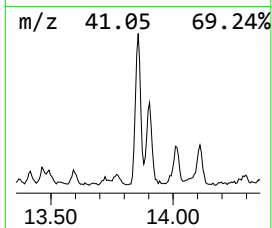
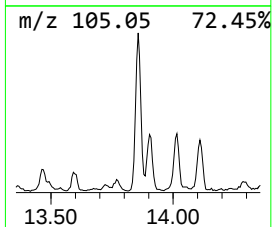
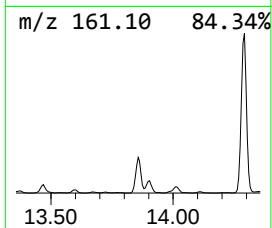
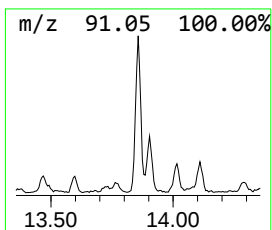
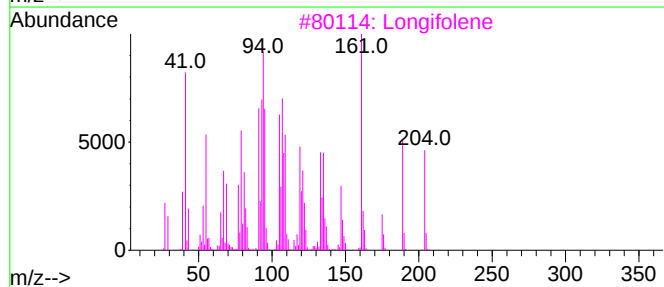
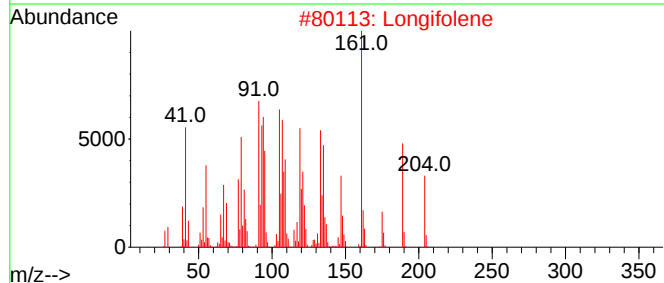
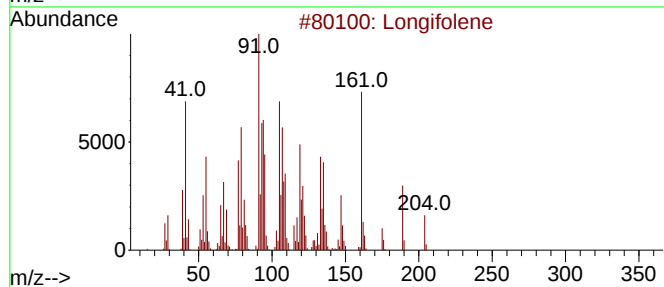
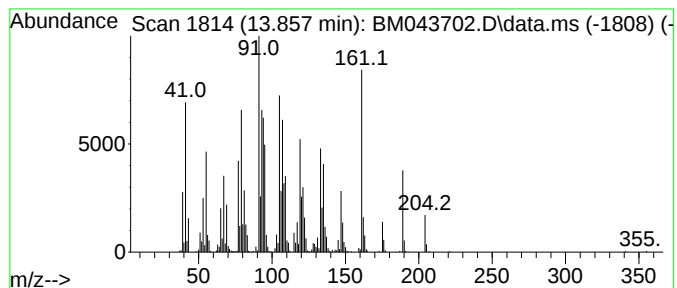
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Quant Title : SVOA CALIBRATION

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

Peak Number 3 Longifolene Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.857	8.83 ng/ul	1069270	Acenaphthene-d10	14.592

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Longifolene	204	C15H24	000475-20-7	99
2			Longifolene	204	C15H24	000475-20-7	99
3			Longifolene	204	C15H24	000475-20-7	99
4			(3R,4aS,5R)-4a,5-Dimethyl-3-(pro...	204	C15H24	024741-64-8	97
5			Naphthalene, 1,2,3,5,6,7,8,8a-oc...	204	C15H24	004630-07-3	97



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM010224\
Data File : BM043702.D
Acq On : 02 Jan 2024 19:54
Operator : MA/JU
Sample : 05920-07
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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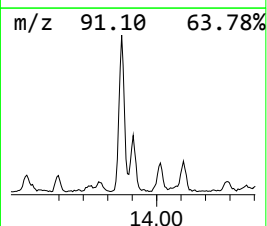
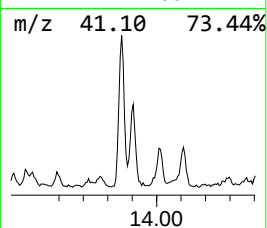
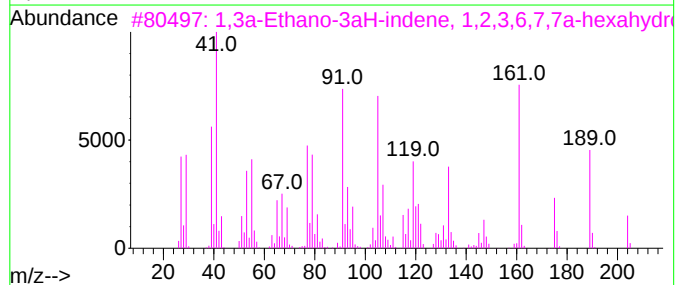
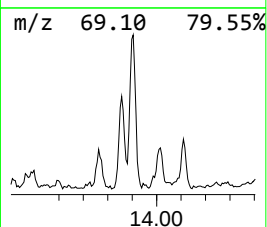
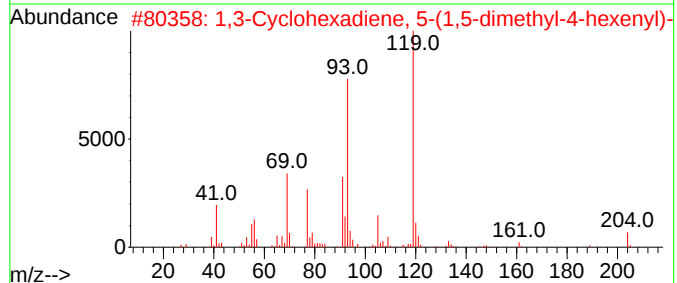
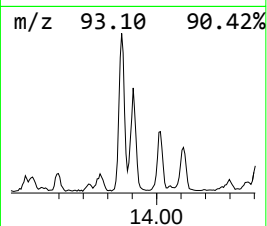
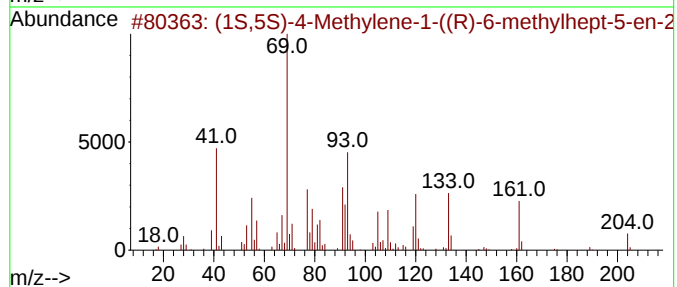
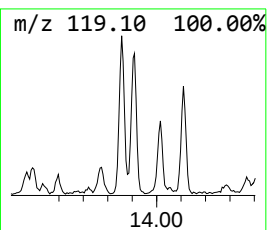
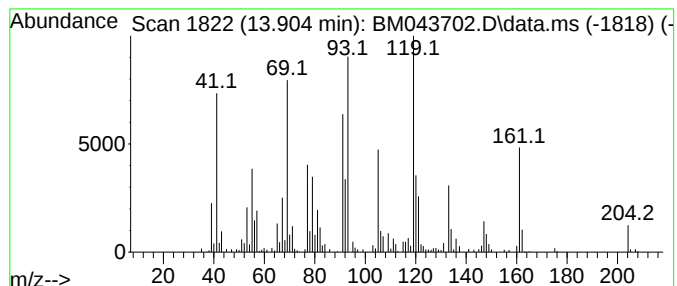
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 (1S,5S)-4-Methylene-1-((R)-... Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.904	3.58 ng/ul	433844	Acenaphthene-d10	14.592

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			(1S,5S)-4-Methylene-1-((R)-6-met...	204	C15H24	058319-04-3	50
2			1,3-Cyclohexadiene, 5-(1,5-dimet...	204	C15H24	000495-60-3	43
3			1,3a-Ethano-3aH-indene, 1,2,3,6,...	204	C15H24	004545-68-0	38
4			(R)-1-Methyl-4-(6-methylhept-5-e...	204	C15H24	028976-67-2	38
5			1H-3a,7-Methanoazulene, 2,3,4,7,...	204	C15H24	000469-61-4	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM010224\
Data File : BM043702.D
Acq On : 02 Jan 2024 19:54
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Sample : 05920-07
Misc :
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Instrument :
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ClientSampleId :
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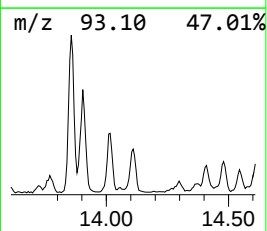
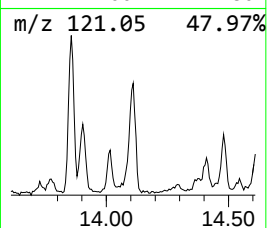
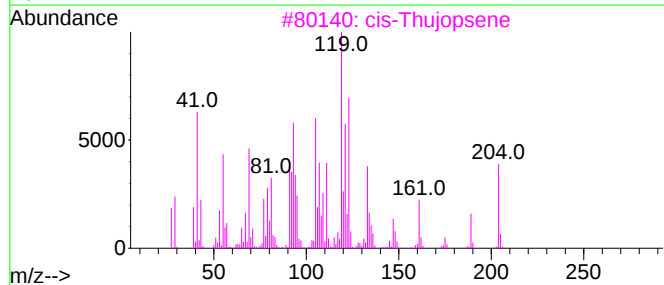
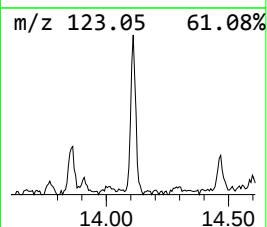
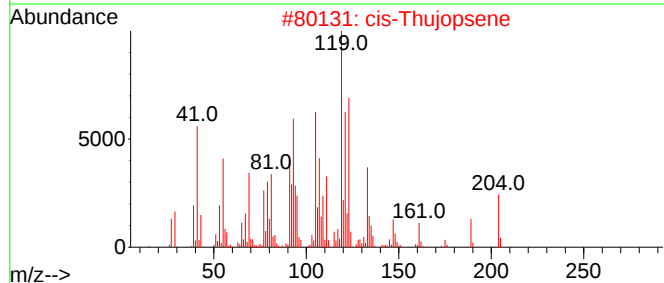
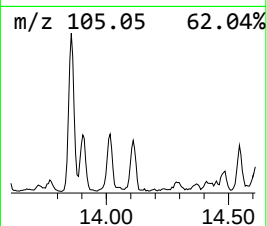
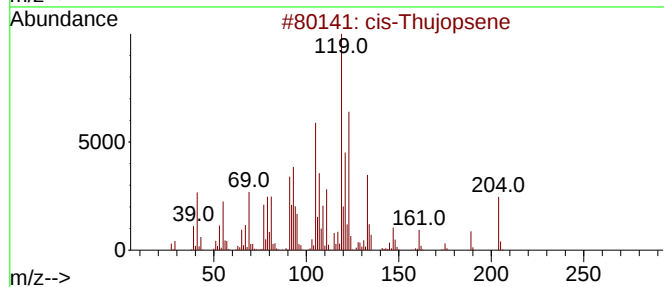
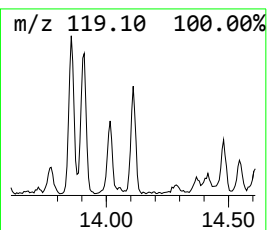
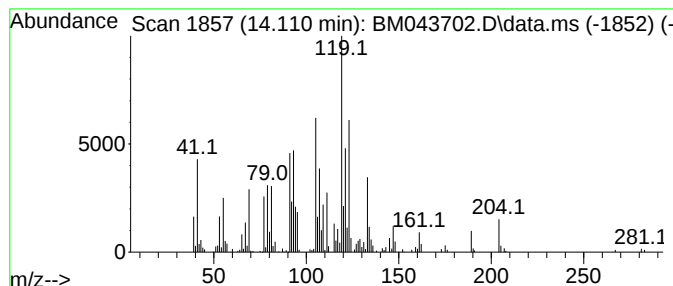
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 cis-Thujopsene Concentration Rank 21

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.110	2.43 ng/ul	293975	Acenaphthene-d10	14.592

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			cis-Thujopsene	204	C15H24	000470-40-6	99
2			cis-Thujopsene	204	C15H24	000470-40-6	95
3			cis-Thujopsene	204	C15H24	000470-40-6	94
4			cis-Thujopsene	204	C15H24	000470-40-6	83
5			Tricyclo[5.4.0.0(2,8)]undec-9-en...	204	C15H24	005989-08-2	78



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM010224\
Data File : BM043702.D
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Sample : 05920-07
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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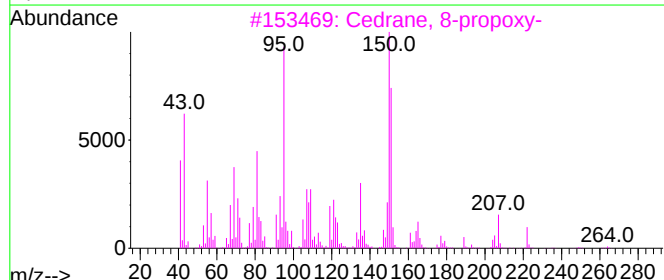
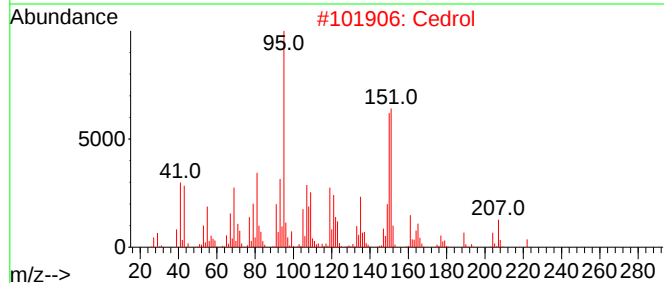
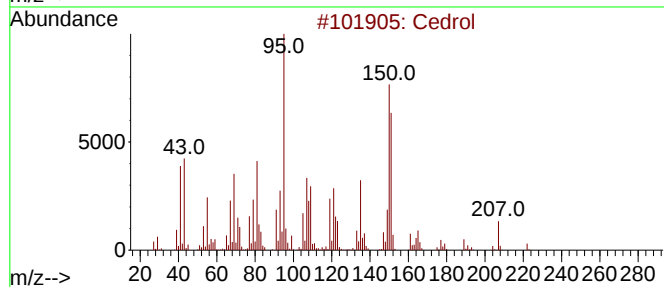
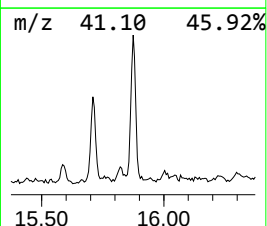
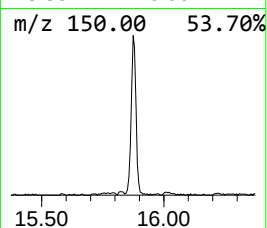
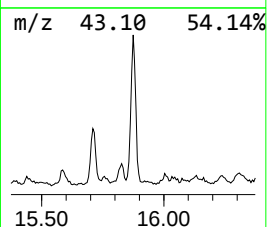
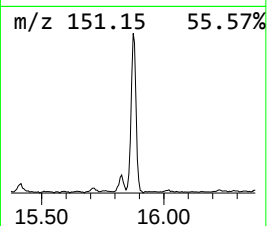
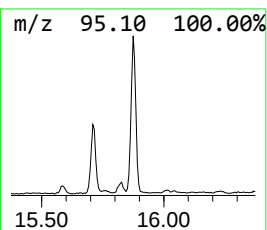
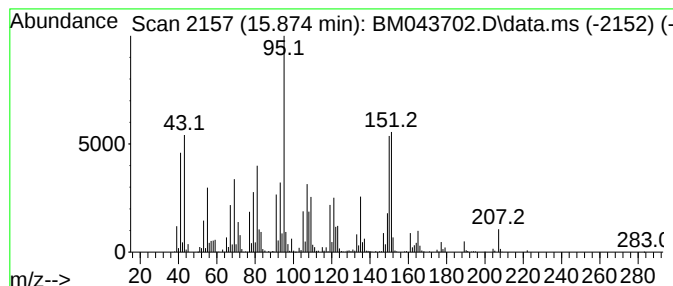
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 Cedrol Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.874	6.82 ng/ul	826501	Acenaphthene-d10	14.592

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cedrol	222	C15H26O	000077-53-2	91
2			Cedrol	222	C15H26O	000077-53-2	91
3			Cedrane, 8-propoxy-	264	C18H32O	019870-75-8	90
4			Cedrol	222	C15H26O	000077-53-2	86
5			Cedrol	222	C15H26O	000077-53-2	83



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM010224\
Data File : BM043702.D
Acq On : 02 Jan 2024 19:54
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Sample : 05920-07
Misc :
ALS Vial : 15 Sample Multiplier: 1

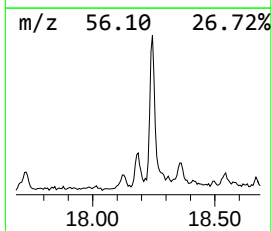
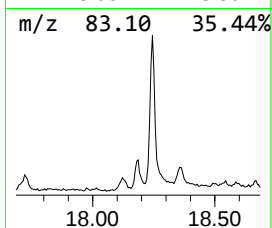
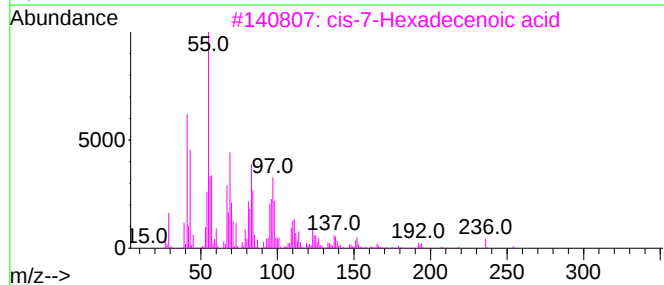
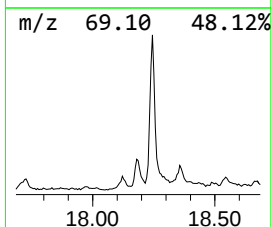
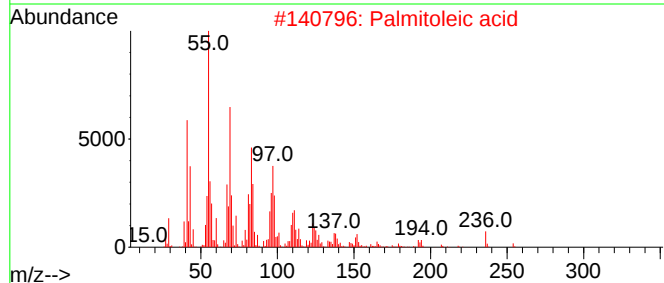
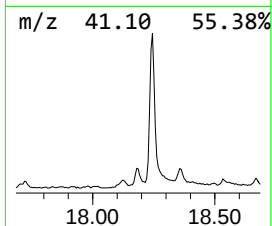
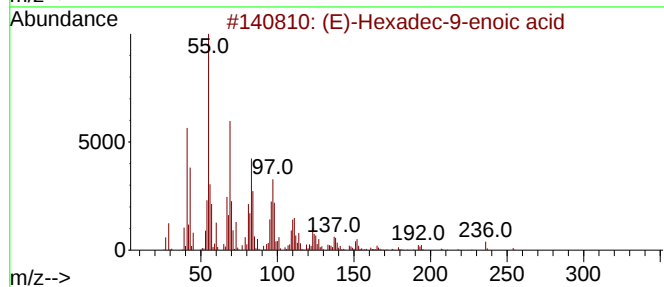
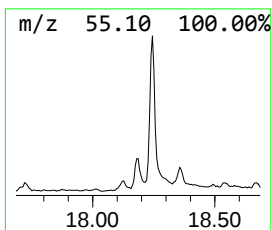
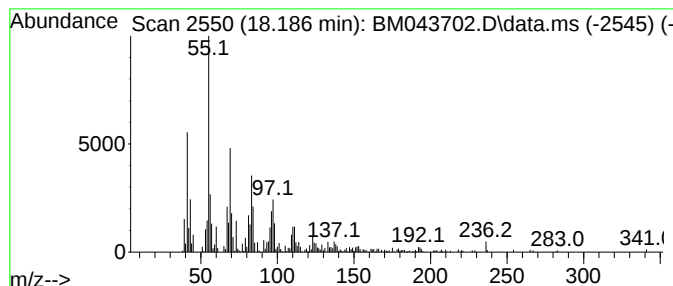
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ClientSampleId :
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 (E)-Hexadec-9-enoic acid Concentration Rank 23

R.T.	EstConc	Area	Relative to ISTD		R.T.	
18.186	2.04 ng/ul	271354	Phenanthrene-d10		17.339	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	(E)-Hexadec-9-enoic acid		254	C16H30O2	010030-73-6	99
2	Palmitoleic acid		254	C16H30O2	000373-49-9	99
3	cis-7-Hexadecenoic acid		254	C16H30O2	002416-19-5	98
4	Hexadecenoic acid, Z-11-		254	C16H30O2	002416-20-8	93
5	9-Hexadecenoic acid		254	C16H30O2	002091-29-4	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM010224\
Data File : BM043702.D
Acq On : 02 Jan 2024 19:54
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Sample : 05920-07
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ALS Vial : 15 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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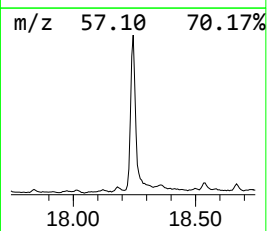
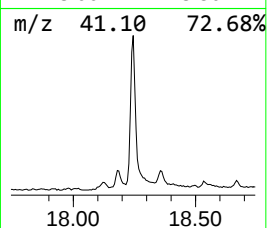
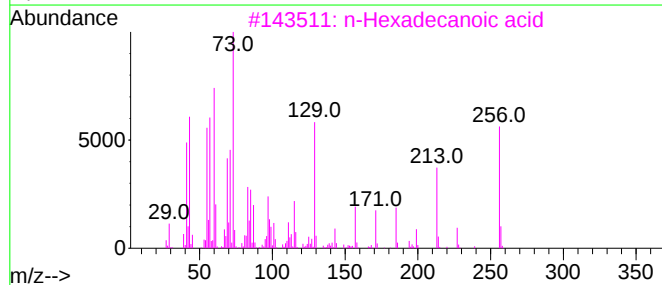
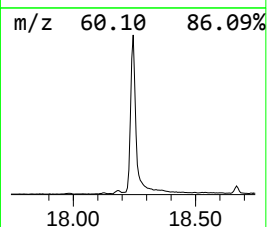
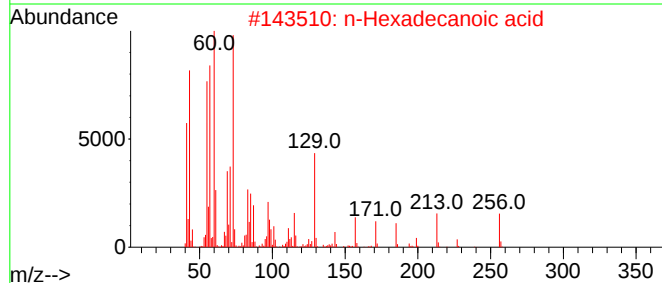
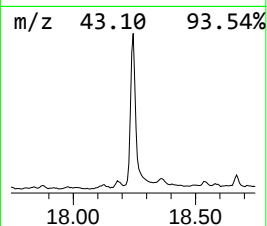
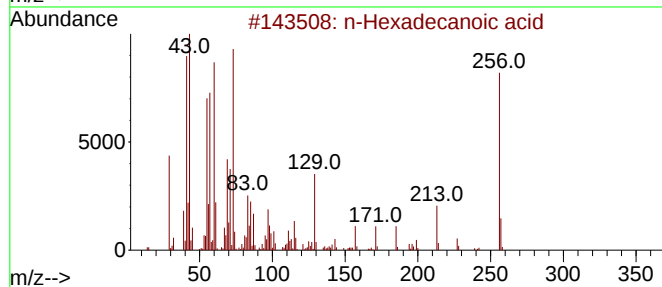
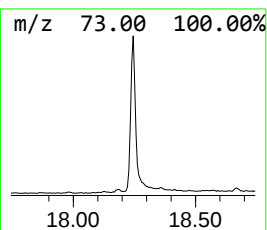
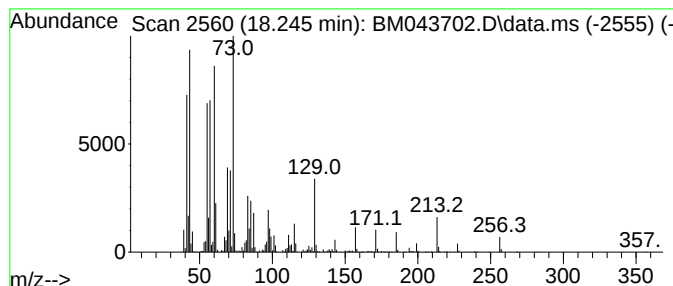
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 n-Hexadecanoic acid Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.245	17.72 ng/ul	2358550	Phenanthrene-d10	17.339

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	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	91		



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM010224\
Data File : BM043702.D
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Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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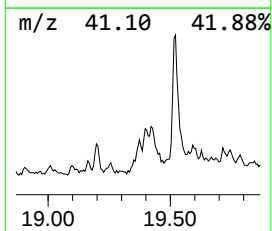
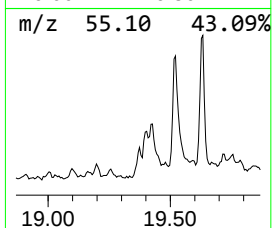
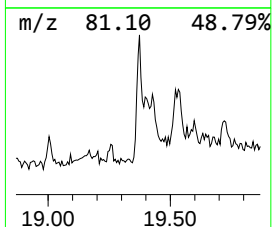
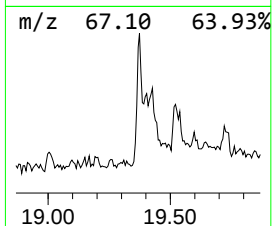
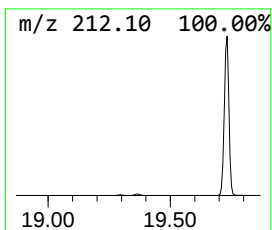
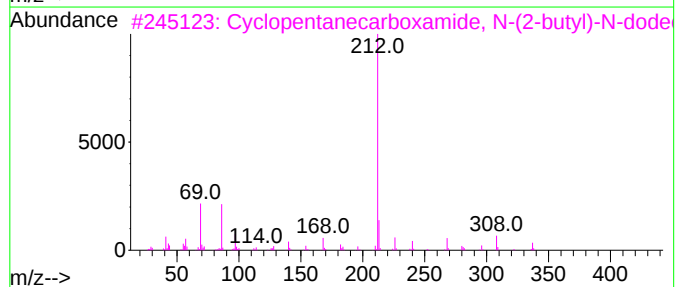
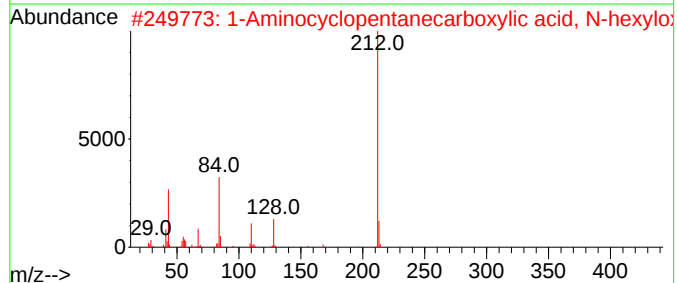
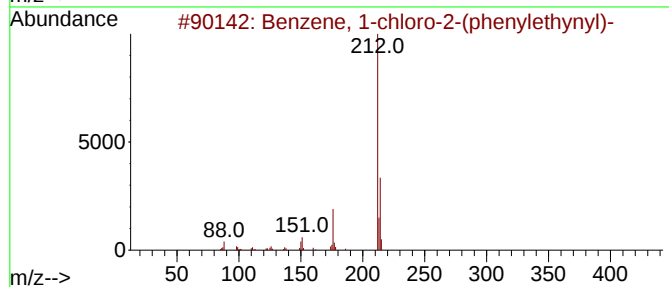
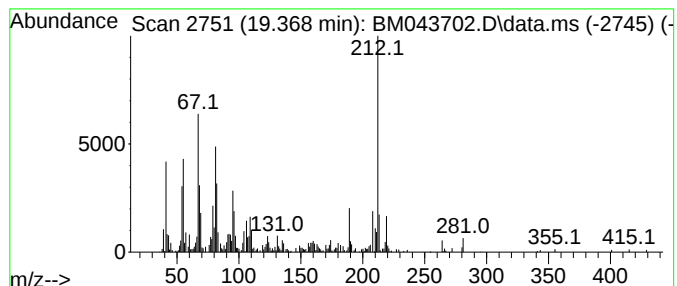
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 unknown-01 Concentration Rank 22

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.368	2.40 ng/ul	319240	Phenanthrene-d10	17.339

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-chloro-2-(phenylethyn...	212	C14H9Cl	010271-57-5	38
2			1-Aminocyclopentanecarboxylic ac...	341	C19H35NO4	1000392-42-8	35
3			Cyclopentanecarboxamide, N-(2-bu...	337	C22H43NO	1000458-93-0	35
4			1-Chloromethyl-3-spiro-cyclohexy...	247	C15H18ClN	1000301-43-2	35
5			1-Aminocyclopentanecarboxylic ac...	383	C22H41NO4	1000392-43-1	35



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM010224\
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ALS Vial : 15 Sample Multiplier: 1

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ClientSampleId :
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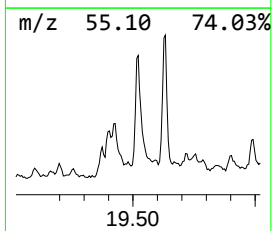
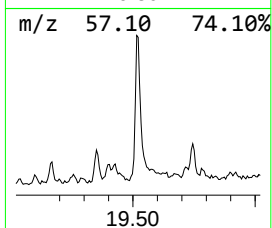
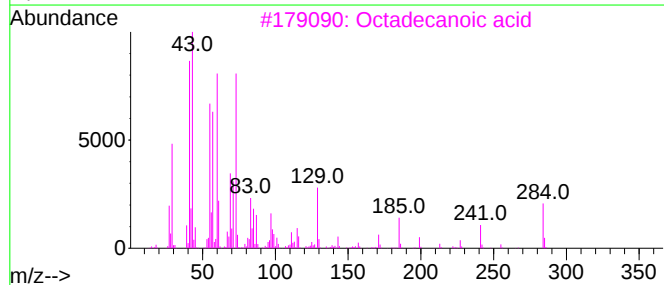
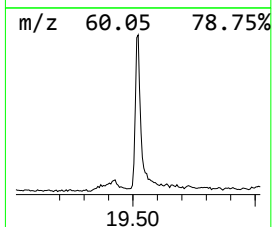
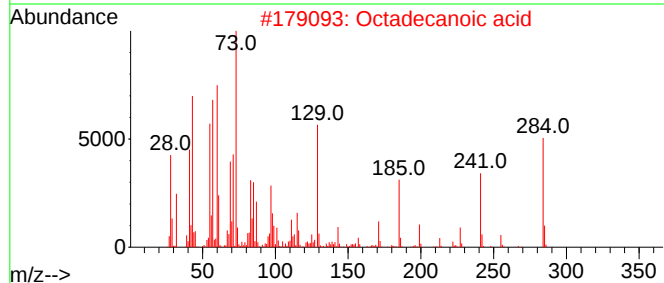
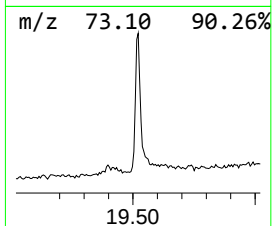
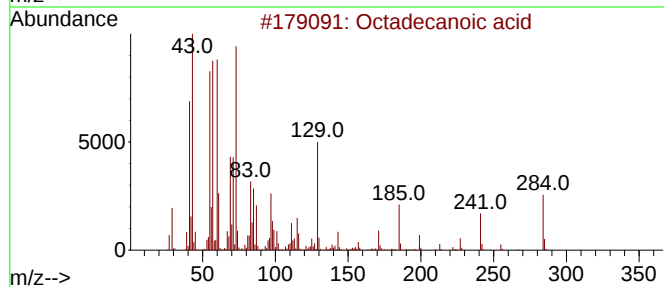
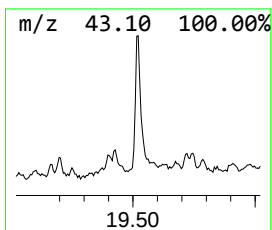
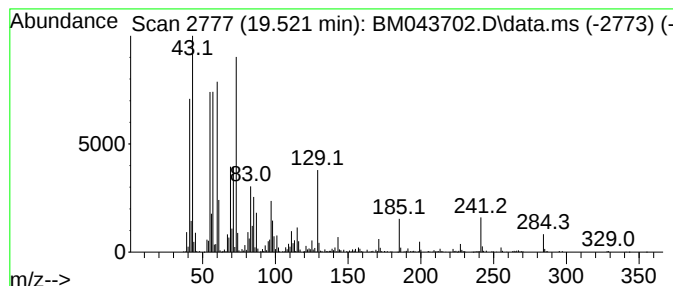
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 Octadecanoic acid Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.521	5.67 ng/ul	817450	Chrysene-d12	21.521

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octadecanoic acid	284	C18H36O2	000057-11-4	99
2			Octadecanoic acid	284	C18H36O2	000057-11-4	99
3			Octadecanoic acid	284	C18H36O2	000057-11-4	99
4			Octadecanoic acid	284	C18H36O2	000057-11-4	98
5			Pentadecanoic acid	242	C15H30O2	001002-84-2	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM010224\
Data File : BM043702.D
Acq On : 02 Jan 2024 19:54
Operator : MA/JU
Sample : 05920-07
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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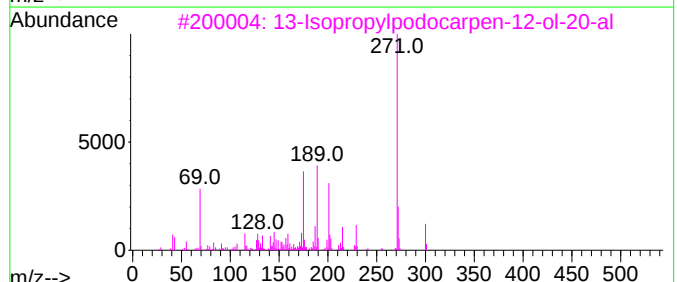
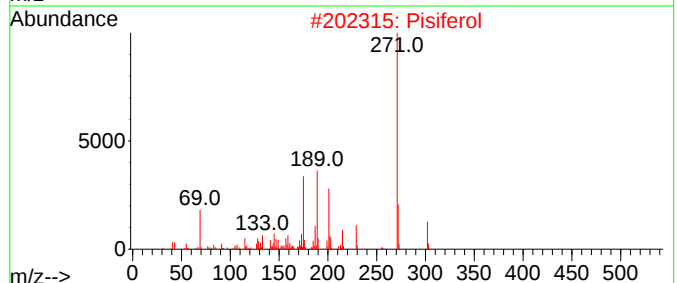
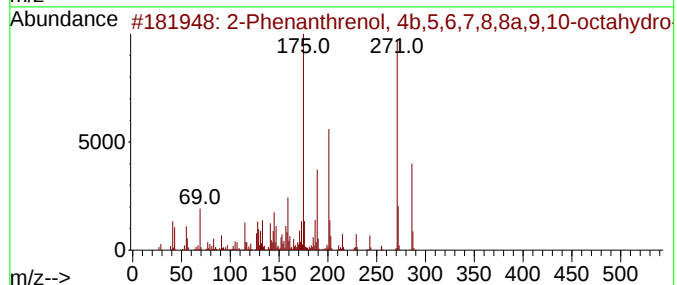
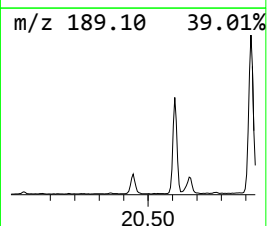
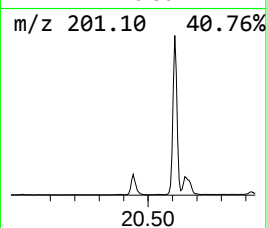
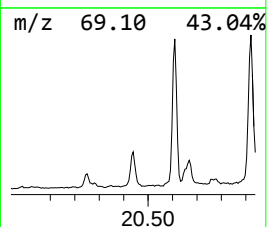
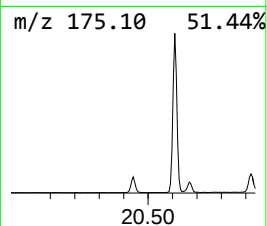
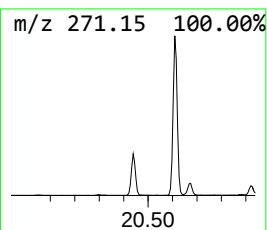
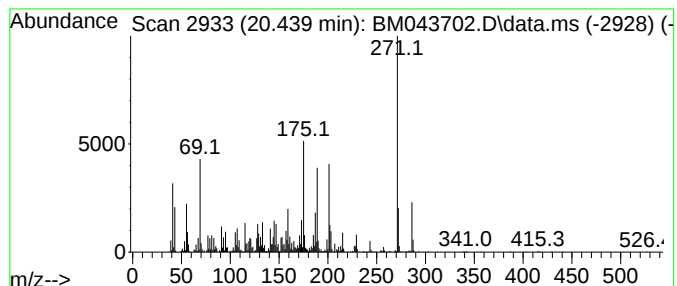
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Quant Title : SVOA CALIBRATION

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

Peak Number 11 2-Phenanthrenol, 4b,5,6,7,8... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.439	7.34 ng/ul	1058840	Chrysene-d12	21.521

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Phenanthrenol, 4b,5,6,7,8,8a,9...	286	C20H30O	000511-15-9	96		
2	Pisiferol	302	C20H30O2	024035-36-7	64		
3	13-Isopropylpodocarpin-12-ol-20-al	300	C20H28O2	024035-37-8	52		
4	13-Isopropylpodocarpin-12-ol-20-al	300	C20H28O2	024035-37-8	49		
5	2-Phenanthrenol, 4b,5,6,7,8,8a,9...	286	C20H30O	000511-15-9	46		



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM010224\
Data File : BM043702.D
Acq On : 02 Jan 2024 19:54
Operator : MA/JU
Sample : 05920-07
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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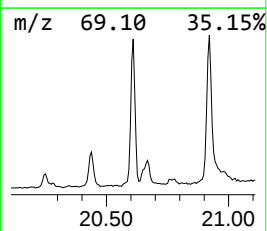
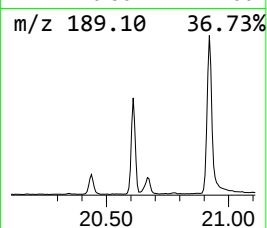
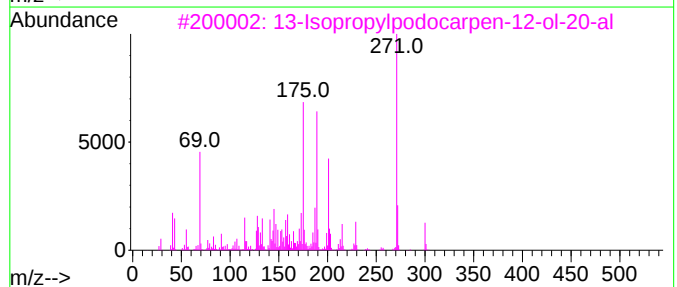
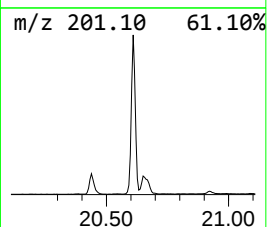
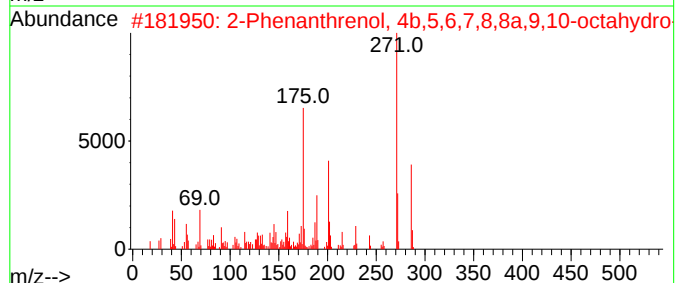
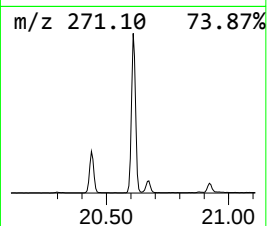
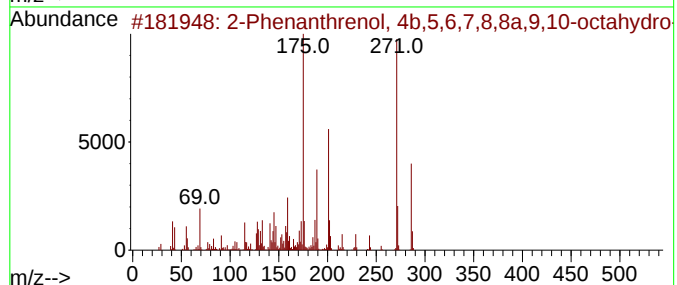
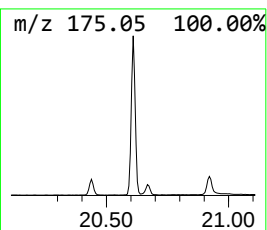
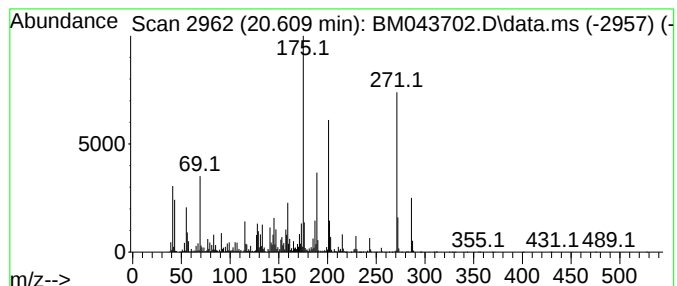
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Quant Title : SVOA CALIBRATION

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

Peak Number 12 13-Isopropylpodocarpin-12-o... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.609	26.67 ng/ul	3845470	Chrysene-d12	21.521

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Phenanthrenol, 4b,5,6,7,8,8a,9...	286	C20H30O	000511-15-9	99		
2	2-Phenanthrenol, 4b,5,6,7,8,8a,9...	286	C20H30O	000511-15-9	83		
3	13-Isopropylpodocarpin-12-ol-20-al	300	C20H28O2	024035-37-8	55		
4	Oxazole, 2-(1-naphthalenyl)-5-ph...	271	C19H13NO	000846-63-9	35		
5	(-)-Morphinan-6-one, 4-methoxy-	271	C17H21NO2	1000129-09-1	20		



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM010224\
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Operator : MA/JU
Sample : 05920-07
Misc :
ALS Vial : 15 Sample Multiplier: 1

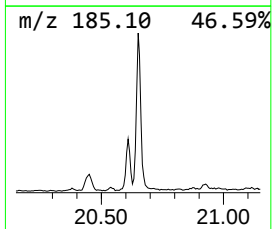
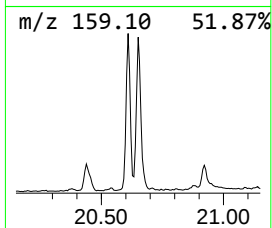
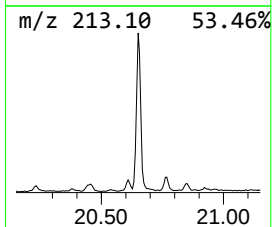
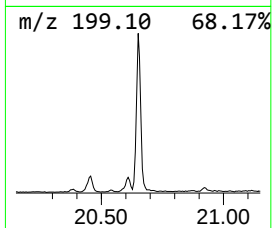
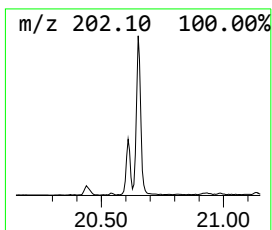
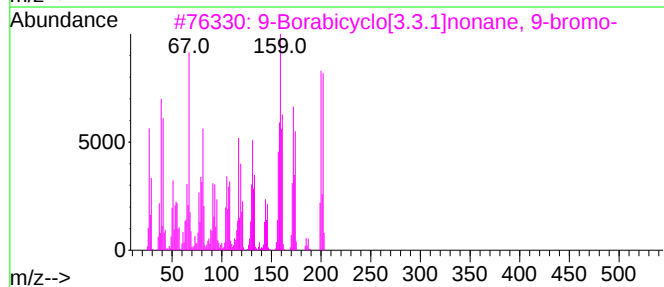
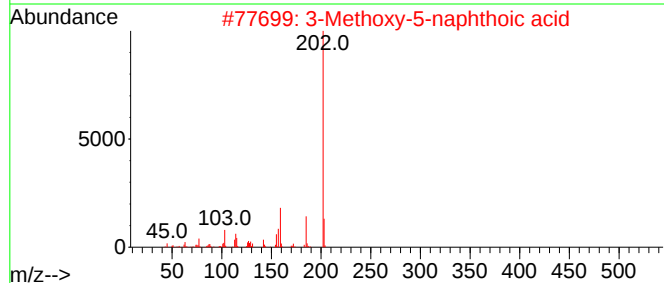
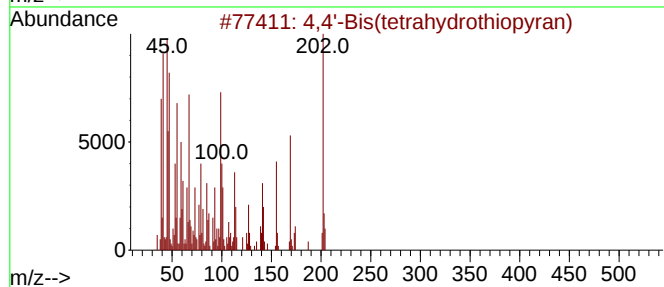
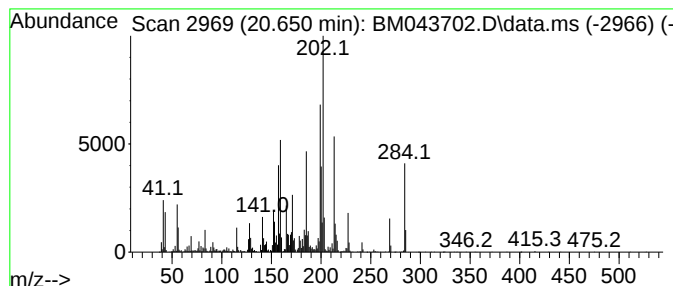
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ClientSampleId :
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Quant Title : SVOA CALIBRATION

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

Peak Number 13 4,4'-Bis(tetrahydrothiopyran) Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.	
20.650	15.58 ng/ul	2246060	Chrysene-d12	21.521	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4,4'-Bis(tetrahydrothiopyran)	202	C10H18S2	116196-83-9	91
2	3-Methoxy-5-naphthoic acid	202	C12H10O3	007498-58-0	46
3	9-Borabicyclo[3.3.1]nonane, 9-br...	200	C8H14BBr	022086-45-9	38
4	5-Hydroxy-3-methyl-1-phenylpyraz...	202	C11H10N2O2	060484-29-9	35
5	[1,2,5]-Oxadiazolo[3,4-a]cinnoli...	202	C10H10N4O	1000260-59-4	20



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM010224\
Data File : BM043702.D
Acq On : 02 Jan 2024 19:54
Operator : MA/JU
Sample : 05920-07
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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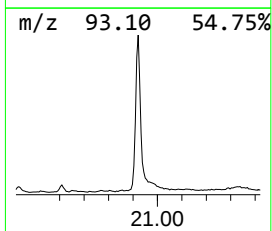
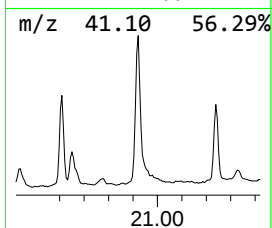
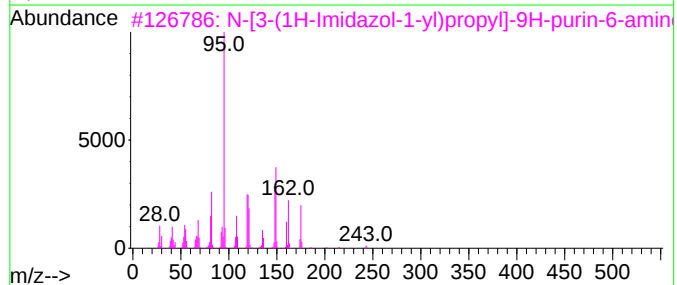
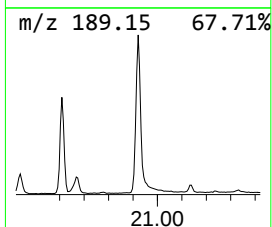
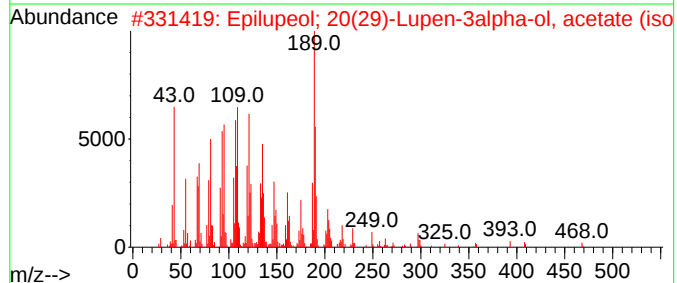
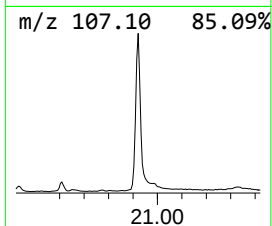
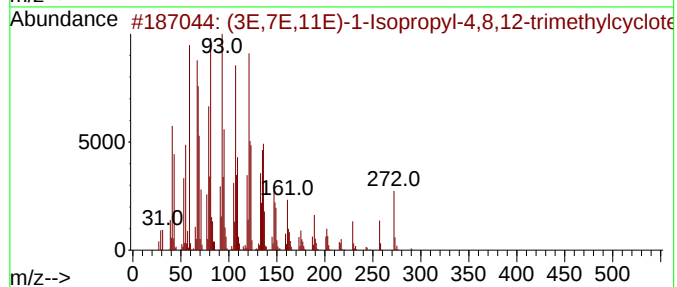
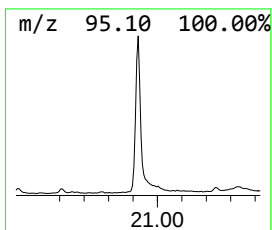
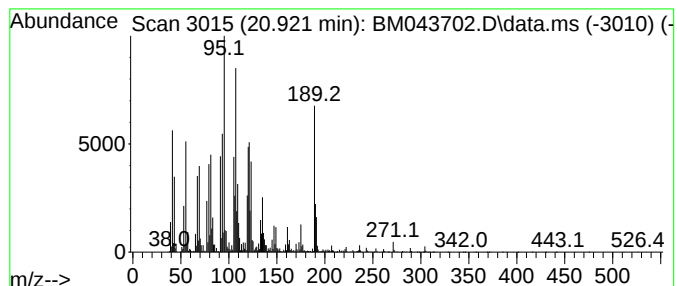
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Quant Title : SVOA CALIBRATION

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

Peak Number 14 unknown-02 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.921	50.77 ng/ul	7320400	Chrysene-d12	21.521

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			(3E,7E,11E)-1-Isopropyl-4,8,12-t...	290	C20H34O	067921-02-2	43
2			Epilupeol; 20(29)-Lupen-3alpha-o...	468	C32H52O2	1010513-01-8	30
3			N-[3-(1H-Imidazol-1-yl)propyl]-9...	243	C11H13N7	537666-87-8	25
4			Cyclopentene, 1-isopropyl-2,3-di...	138	C10H18	007712-73-4	14
5			2-Trichloromethyl-1-bicyclo[2.2....	212	C8H11Cl3	090086-53-6	11



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM010224\
Data File : BM043702.D
Acq On : 02 Jan 2024 19:54
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Sample : 05920-07
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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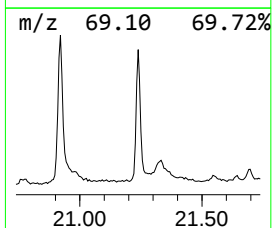
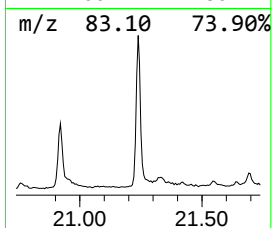
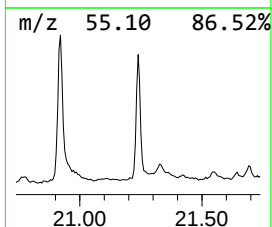
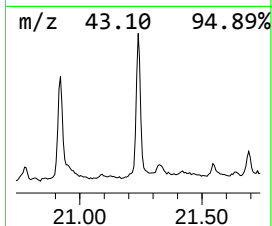
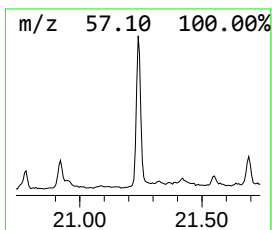
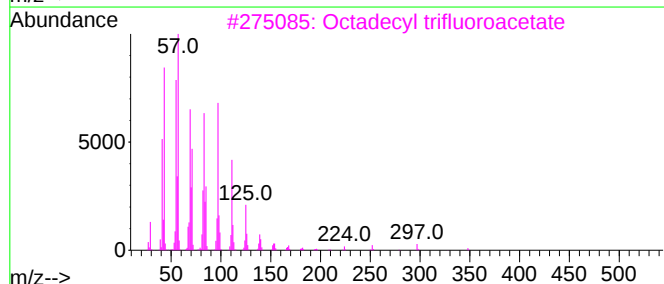
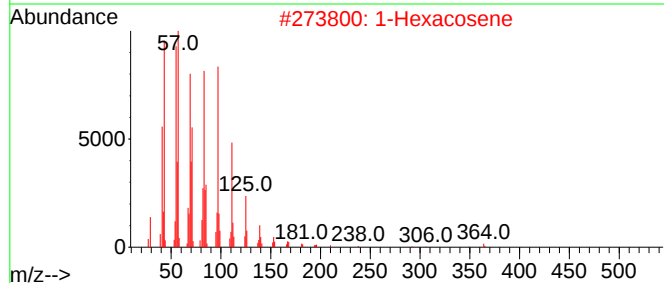
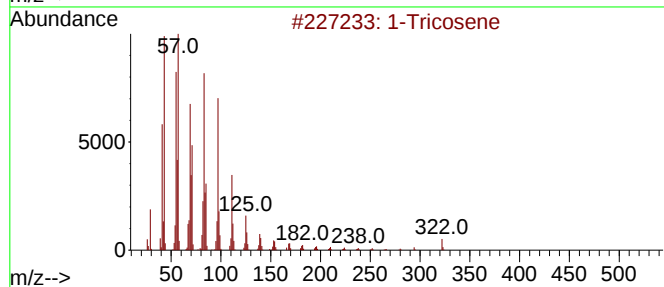
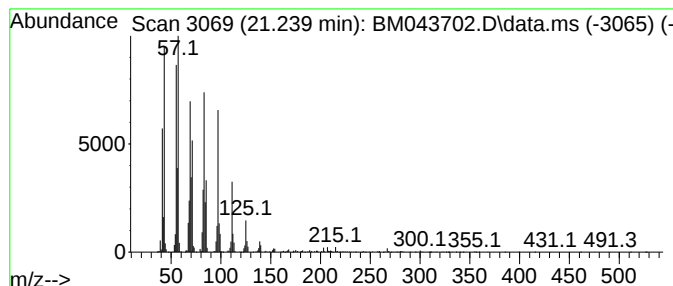
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Quant Title : SVOA CALIBRATION

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

Peak Number 15 (DEL) Alkane: Straight-Chai... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.239	12.22 ng/ul	1761430	Chrysene-d12	21.521

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		1-Tricosene	322	C23H46	018835-32-0	95
2		1-Hexacosene	364	C26H52	018835-33-1	94
3		Octadecyl trifluoroacetate	366	C20H37F3O2	079392-43-1	94
4		Pentadecafluorooctanoic acid, oc...	666	C26H37F15O2	1000406-04-8	93
5		Cyclopentadecane	210	C15H30	000295-48-7	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM010224\
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Sample : 05920-07
Misc :
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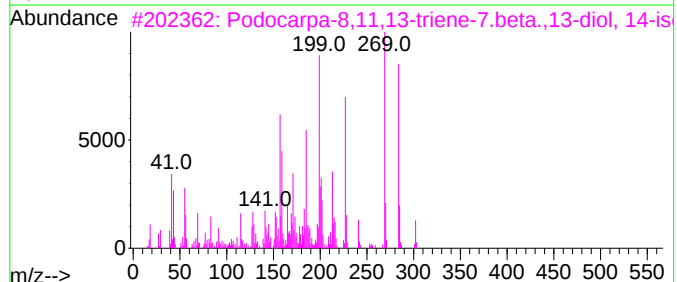
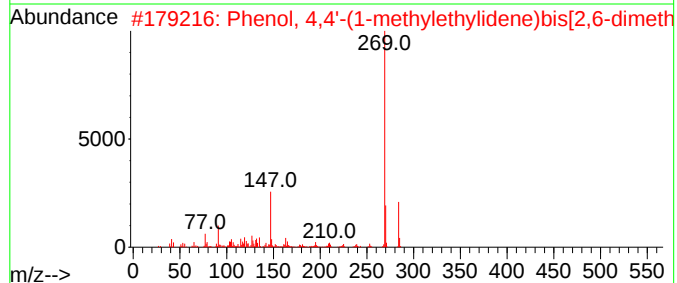
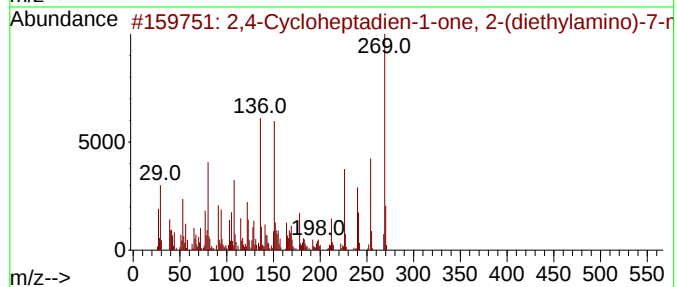
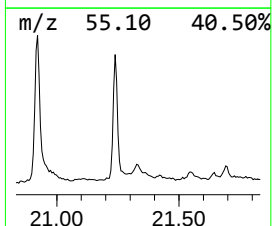
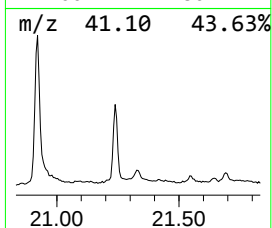
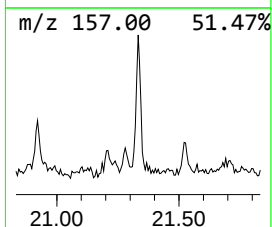
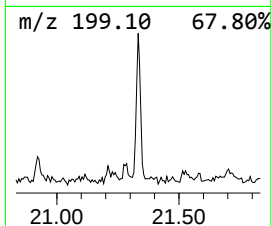
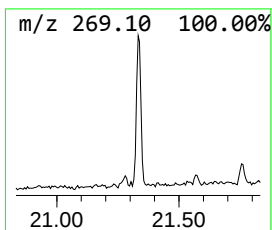
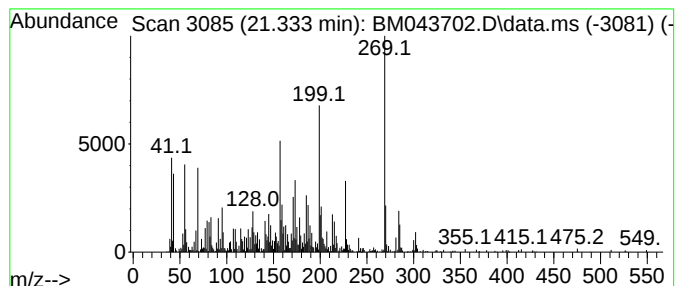
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Quant Title : SVOA CALIBRATION

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

Peak Number 16 2,4-Cycloheptadien-1-one, 2... Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD		R.T.	
21.333	5.83 ng/ul	839993	Chrysene-d12		21.521	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	2,4-Cycloheptadien-1-one, 2-(die...		269	C18H23NO	133436-07-4	78
2	Phenol, 4,4'-(1-methylethylidene...		284	C19H24O2	005613-46-7	60
3	Podocarpa-8,11,13-triene-7.beta....		302	C20H30O2	024338-19-0	43
4	4-Cumylphenol, TMS derivative		284	C18H24OSi	261502-93-6	38
5	Phenol, 4,4'-(1-methylethylidene...		284	C19H24O2	005613-46-7	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM010224\
Data File : BM043702.D
Acq On : 02 Jan 2024 19:54
Operator : MA/JU
Sample : 05920-07
Misc :
ALS Vial : 15 Sample Multiplier: 1

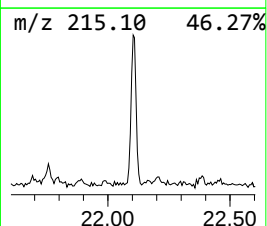
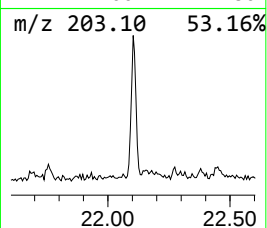
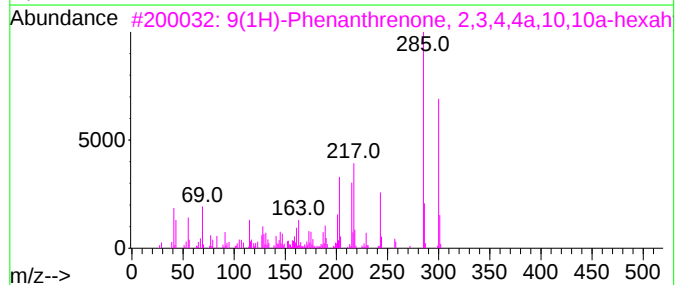
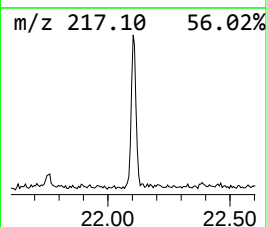
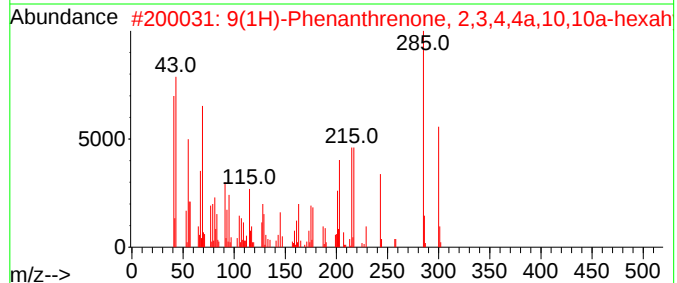
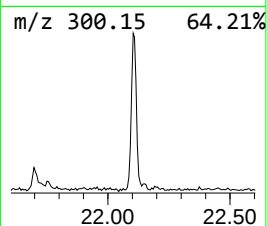
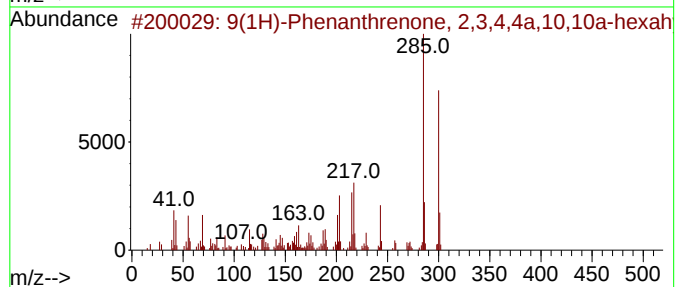
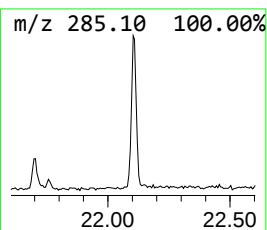
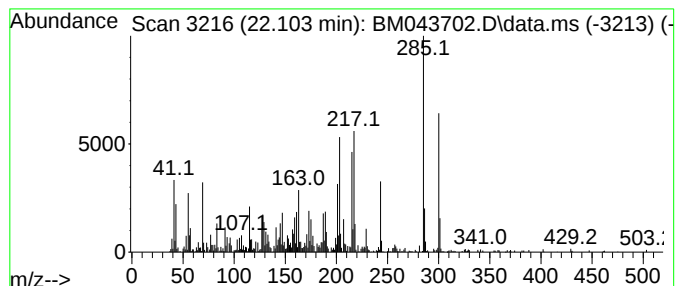
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Quant Title : SVOA CALIBRATION

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

Peak Number 17 9(1H)-Phenanthrenone, 2,3,4... Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.	
22.103	2.61 ng/ul	376319	Chrysene-d12	21.521	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	9(1H)-Phenanthrenone, 2,3,4,4a,1...	300	C20H28O2	000511-05-7	95
2	9(1H)-Phenanthrenone, 2,3,4,4a,1...	300	C20H28O2	000511-05-7	91
3	9(1H)-Phenanthrenone, 2,3,4,4a,1...	300	C20H28O2	000511-05-7	90
4	Phenanthrene, 1,2,3,4,4a,9,10,10...	300	C21H32O	010064-26-3	62
5	Phenanthrene, 1,2,3,4,4a,9,10,10...	300	C21H32O	010064-26-3	58



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM010224\
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ClientSampleId :
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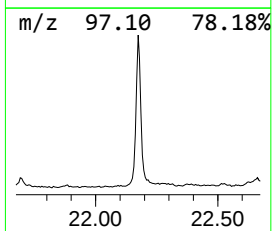
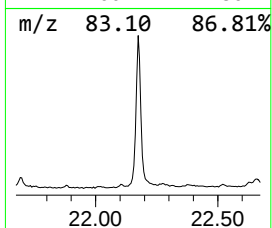
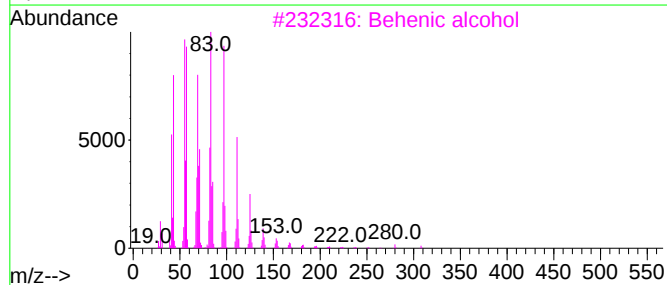
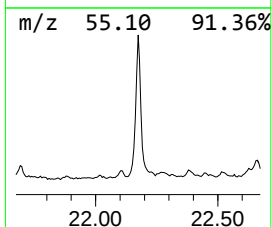
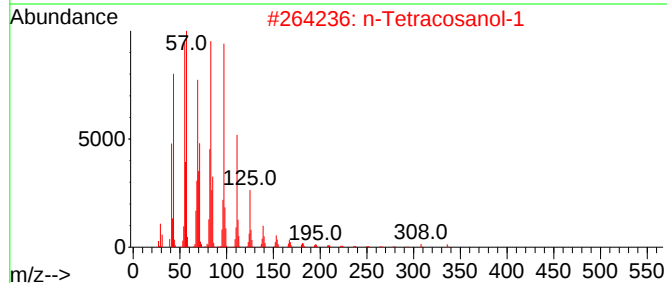
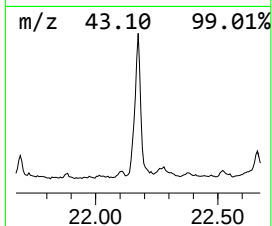
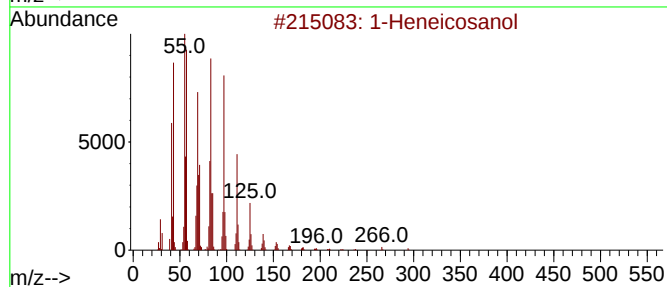
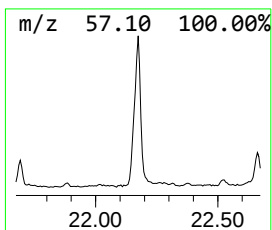
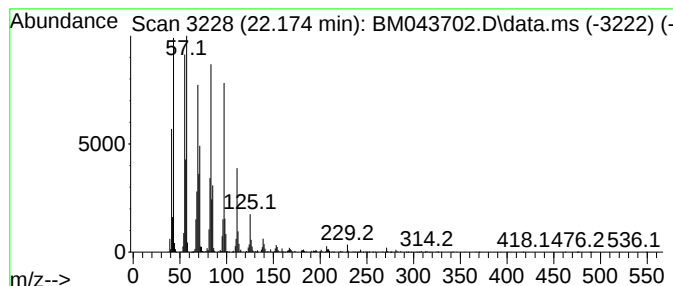
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Quant Title : SVOA CALIBRATION

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

Peak Number 18 1-Heneicosanol Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.174	20.82 ng/ul	3001240	Chrysene-d12	21.521

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		1-Heneicosanol	312	C21H44O	015594-90-8	95
2		n-Tetracosanol-1	354	C24H50O	000506-51-4	94
3		Behenic alcohol	326	C22H46O	000661-19-8	93
4		3-Eicosene, (E)-	280	C20H40	074685-33-9	91
5		1-Hexacosene	364	C26H52	018835-33-1	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM010224\
Data File : BM043702.D
Acq On : 02 Jan 2024 19:54
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Sample : 05920-07
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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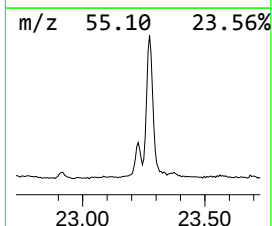
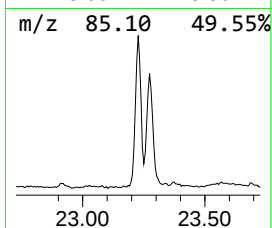
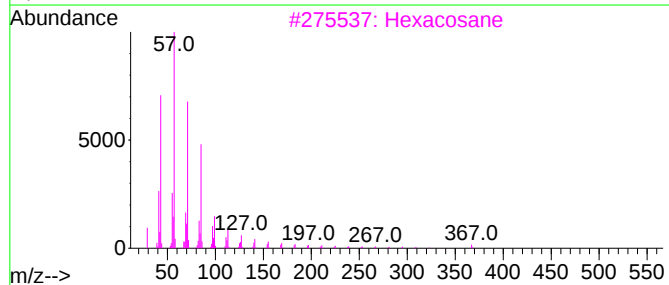
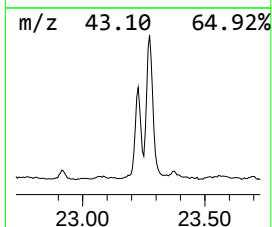
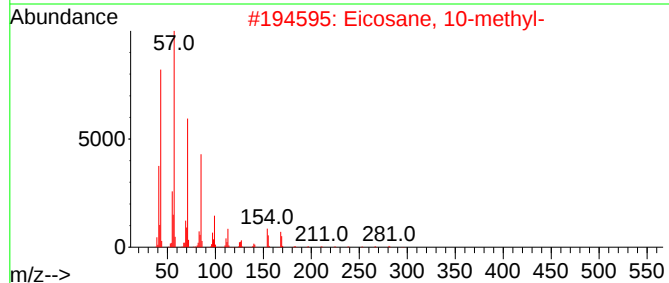
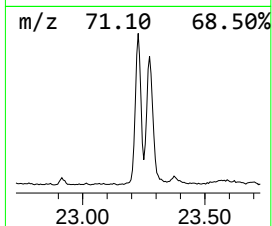
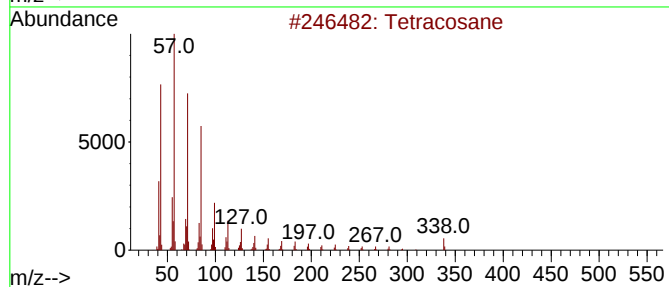
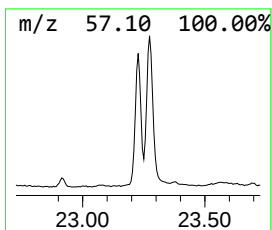
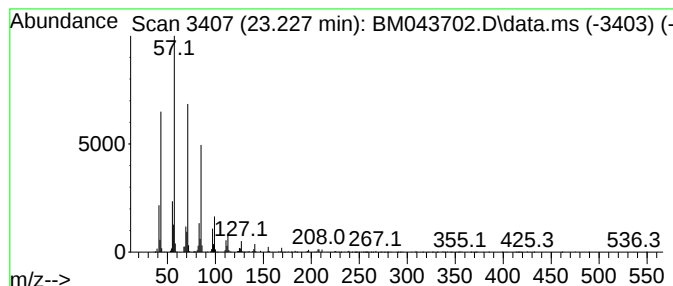
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Quant Title : SVOA CALIBRATION

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

Peak Number 19 (DEL) Alkane: Straight-Chai... Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.227	6.83 ng/ul	1029720	Perylene-d12	23.938

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Tetracosane	338	C24H50	000646-31-1	97
2		Eicosane, 10-methyl-	296	C21H44	054833-23-7	91
3		Hexacosane	366	C26H54	000630-01-3	91
4		Tetracosane	338	C24H50	000646-31-1	90
5		Heneicosane	296	C21H44	000629-94-7	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM010224\
Data File : BM043702.D
Acq On : 02 Jan 2024 19:54
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Sample : 05920-07
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
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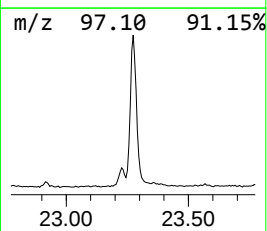
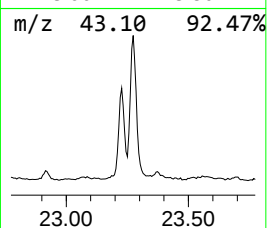
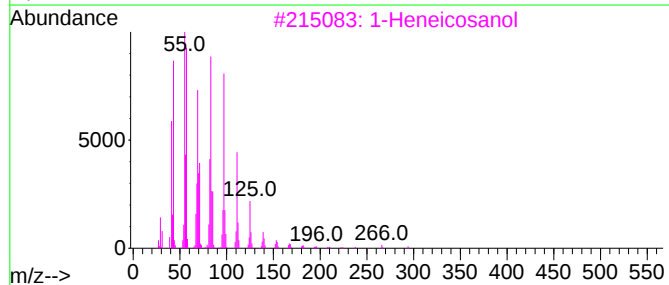
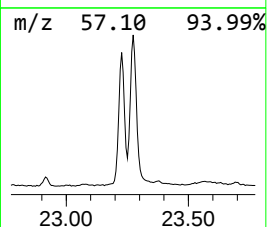
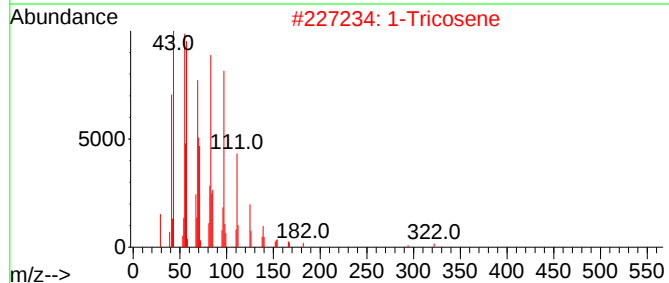
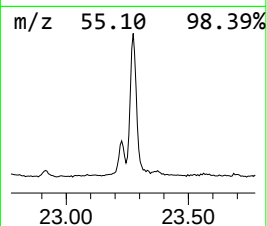
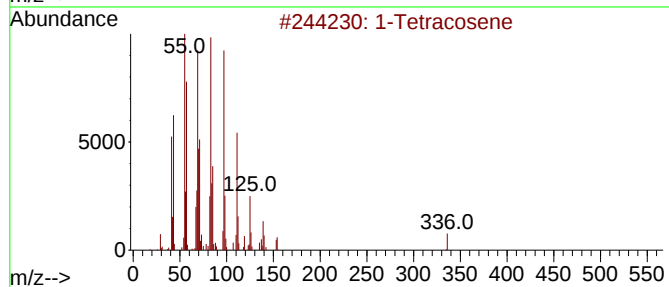
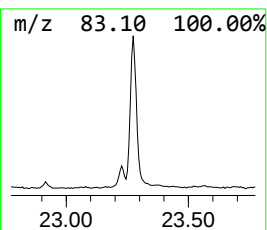
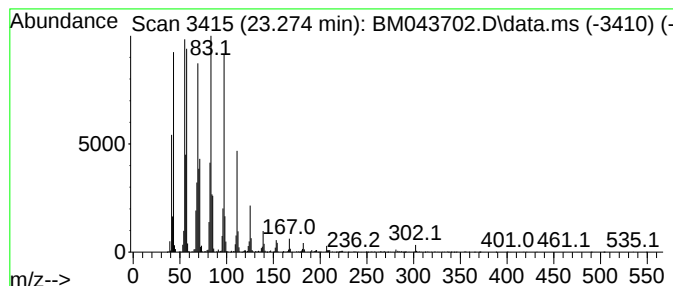
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Quant Title : SVOA CALIBRATION

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

Peak Number 20 (DEL) Alkane: Straight-Chai... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.274	23.18 ng/ul	3496590	Perylene-d12	23.938

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		1-Tetracosene	336	C24H48	010192-32-2	98
2		1-Tricosene	322	C23H46	018835-32-0	95
3		1-Heneicosanol	312	C21H44O	015594-90-8	95
4		n-Tetracosanol-1	354	C24H50O	000506-51-4	94
5		1-Heptacosanol	396	C27H56O	002004-39-9	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM010224\
Data File : BM043702.D
Acq On : 02 Jan 2024 19:54
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Sample : 05920-07
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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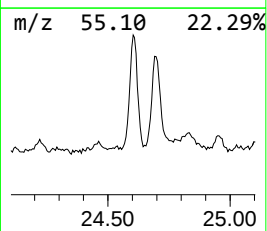
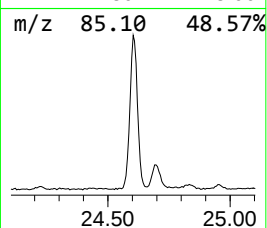
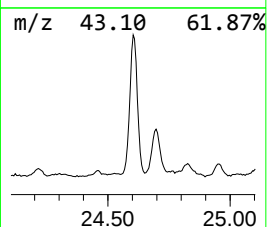
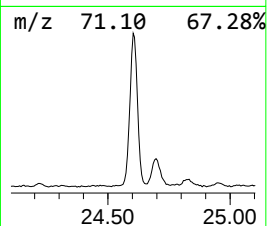
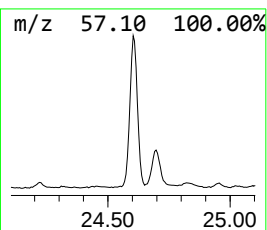
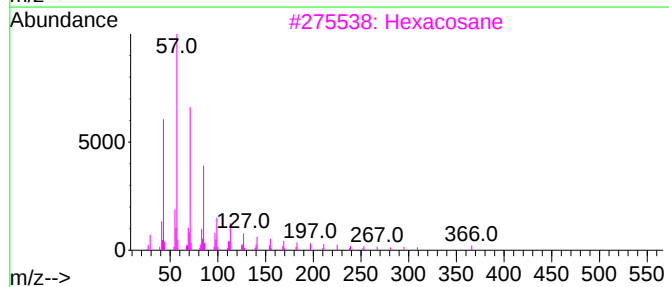
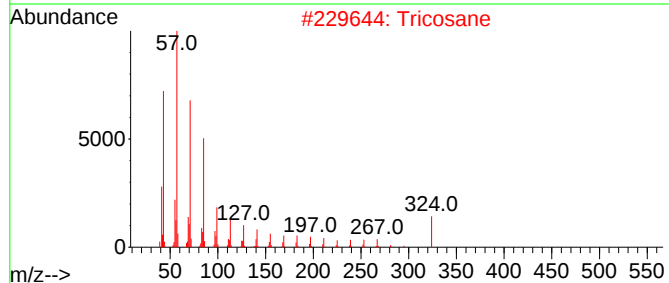
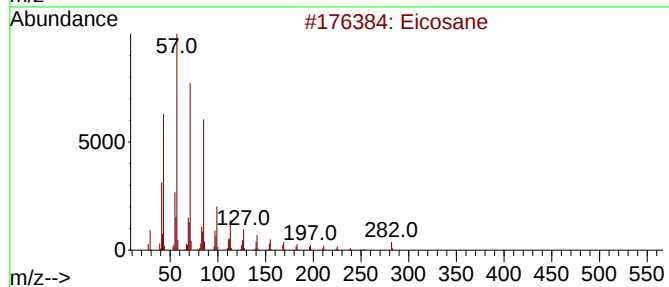
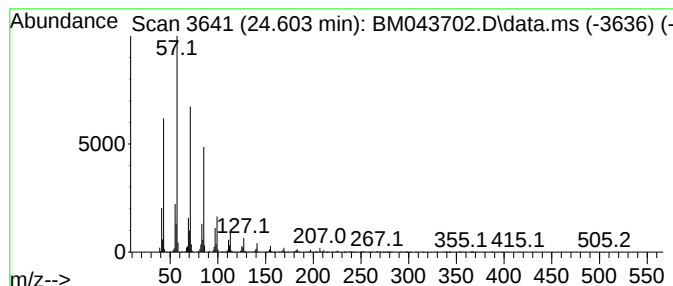
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Quant Title : SVOA CALIBRATION

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

Peak Number 21 (DEL) Alkane: Straight-Chai... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.603	15.13 ng/ul	2282510	Perylene-d12	23.938

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Eicosane	282	C20H42	000112-95-8	95
2			Tricosane	324	C23H48	000638-67-5	93
3			Hexacosane	366	C26H54	000630-01-3	91
4			Dotriacontane, 1-iodo-	576	C32H65I	1000406-32-4	91
5			Octacosane, 2-methyl-	408	C29H60	001560-98-1	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM010224\
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Sample : 05920-07
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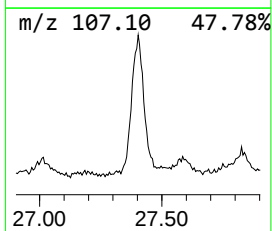
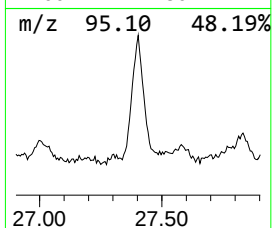
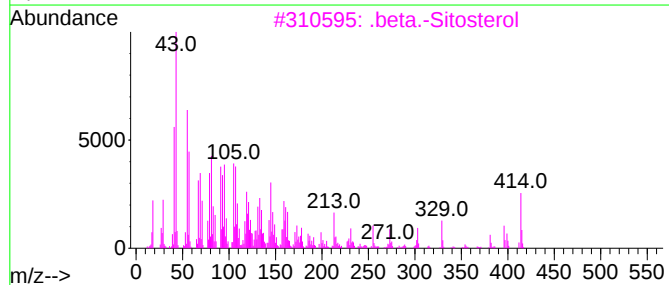
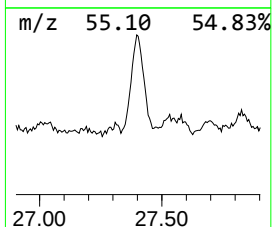
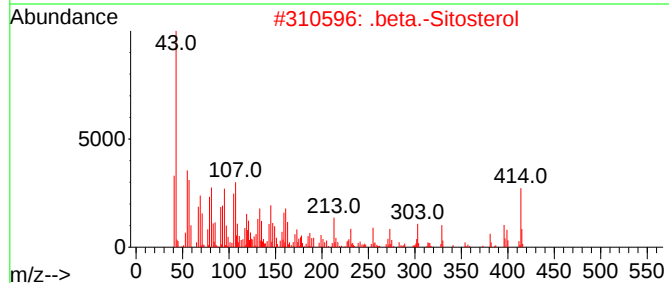
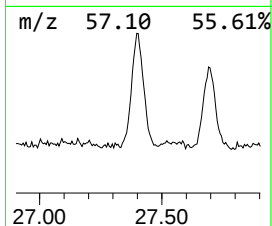
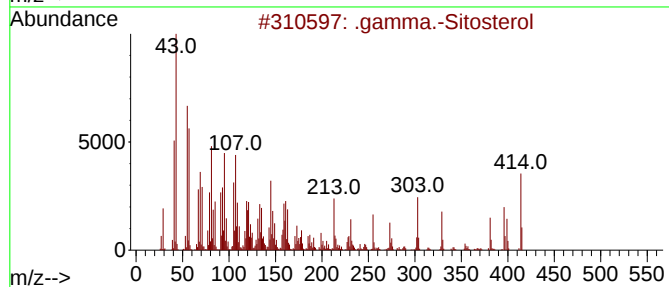
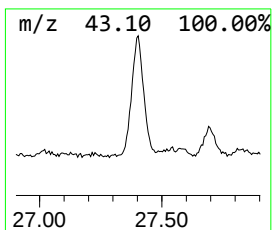
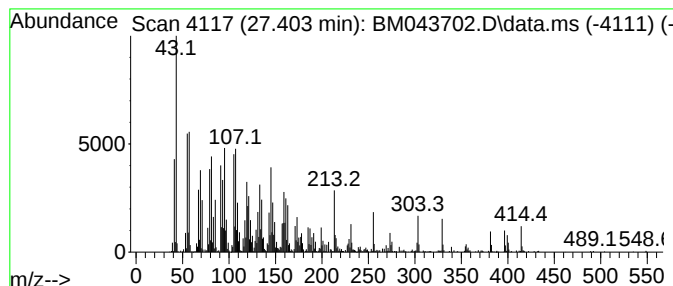
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Quant Title : SVOA CALIBRATION

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

Peak Number 23 .gamma.-Sitosterol Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.	
27.403	19.91 ng/ul	3002530	Perylene-d12	23.938	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	.gamma.-Sitosterol	414	C29H50O	000083-47-6	97
2	.beta.-Sitosterol	414	C29H50O	000083-46-5	94
3	.beta.-Sitosterol	414	C29H50O	000083-46-5	76
4	Isocholesteryl methyl ether	400	C28H48O	029944-53-4	60
5	Ergost-5-en-3-ol, (3.beta.)-	400	C28H48O	004651-51-8	58



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM010224\
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 Operator : MA/JU
 Sample : 05920-07
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
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 ClientSampleId :
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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\NIST0.L
 TIC Integration Parameters: LSCINT.P

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					#	RT	Resp	Conc
trans-.beta.-Oc...	6.622	5.2	ng/ul	356415	1	7.957	1360490	20.0
D-Limonene	8.169	7.3	ng/ul	498679	1	7.957	1360490	20.0
Longifolene	13.857	8.8	ng/ul	1069270	3	14.592	2422270	20.0
(1S,5S)-4-Methy...	13.904	3.6	ng/ul	433844	3	14.592	2422270	20.0
cis-Thujiopsene	14.110	2.4	ng/ul	293975	3	14.592	2422270	20.0
Cedrol	15.874	6.8	ng/ul	826501	3	14.592	2422270	20.0
(E)-Hexadec-9-e...	18.186	2.0	ng/ul	271354	4	17.339	2662680	20.0
n-Hexadecanoic ...	18.245	17.7	ng/ul	2358550	4	17.339	2662680	20.0
unknown-01	19.368	2.4	ng/ul	319240	4	17.339	2662680	20.0
Octadecanoic acid	19.521	5.7	ng/ul	817450	5	21.521	2883590	20.0
2-Phenanthrenol...	20.439	7.3	ng/ul	1058840	5	21.521	2883590	20.0
13-Isopropylpod...	20.609	26.7	ng/ul	3845470	5	21.521	2883590	20.0
4,4'-Bis(tetra...	20.650	15.6	ng/ul	2246060	5	21.521	2883590	20.0
unknown-02	20.921	50.8	ng/ul	7320400	5	21.521	2883590	20.0
(DEL) Alkane: S...	21.239	12.2	ng/ul	1761430	5	21.521	2883590	20.0
2,4-Cycloheptad...	21.333	5.8	ng/ul	839993	5	21.521	2883590	20.0
9(1H)-Phenanthr...	22.103	2.6	ng/ul	376319	5	21.521	2883590	20.0
1-Heneicosanol	22.174	20.8	ng/ul	3001240	5	21.521	2883590	20.0
(DEL) Alkane: S...	23.227	6.8	ng/ul	1029720	6	23.938	3016270	20.0
(DEL) Alkane: S...	23.274	23.2	ng/ul	3496590	6	23.938	3016270	20.0
(DEL) Alkane: S...	24.603	15.1	ng/ul	2282510	6	23.938	3016270	20.0
Triacontan-15-ol	24.697	7.1	ng/ul	1075670	6	23.938	3016270	20.0
.gamma.-Sitosterol	27.403	19.9	ng/ul	3002530	6	23.938	3016270	20.0