

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM010224\
 Data File : BM043748.D
 Acq On : 04 Jan 2024 00:09
 Operator : MA/JU
 Sample : 05951-04
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 GCF68

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: BM043748.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.375	29	32	46	rVB	82701	144893	1.74%	0.111%
2	3.675	80	83	92	rVB	46031	80193	0.96%	0.062%
3	3.805	101	105	118	rBV	701833	1177165	14.14%	0.903%
4	5.311	356	361	370	rBV	128650	236870	2.85%	0.182%
5	6.452	548	555	563	rBV	118914	194464	2.34%	0.149%
6	6.616	574	583	591	rBV	173579	295900	3.56%	0.227%
7	7.140	662	672	682	rBV	1155713	2002151	24.06%	1.536%
8	7.246	682	690	695	rVV	3850095	5927101	71.22%	4.547%
9	7.299	695	699	708	rVV	895731	1413649	16.99%	1.084%
10	7.410	709	718	725	rVB	284833	436557	5.25%	0.335%
11	7.493	725	732	739	rBV	1155734	1843550	22.15%	1.414%
12	7.957	800	811	819	rBV	1175570	1904169	22.88%	1.461%
13	8.163	838	846	852	rBV	156855	268728	3.23%	0.206%
14	8.463	892	897	903	rVB4	20629	35588	0.43%	0.027%
15	8.616	918	923	928	rVB4	35326	54539	0.66%	0.042%
16	8.687	928	935	946	rBV	1211117	2042635	24.54%	1.567%
17	8.846	959	962	972	rVB8	20426	37928	0.46%	0.029%
18	9.134	1005	1011	1024	rVV	1004917	1786930	21.47%	1.371%
19	9.857	1124	1134	1147	rBV	856678	1566385	18.82%	1.202%
20	10.398	1219	1226	1247	rBV	1185682	2299262	27.63%	1.764%
21	10.769	1277	1289	1299	rVV	1533343	2618206	31.46%	2.008%
22	10.928	1309	1316	1334	rBV	308869	737178	8.86%	0.566%
23	11.528	1413	1418	1422	rBV2	28693	52338	0.63%	0.040%
24	11.581	1422	1427	1433	rVB3	52806	94648	1.14%	0.073%
25	12.416	1566	1569	1578	rBV2	19715	45770	0.55%	0.035%
26	12.516	1580	1586	1594	rBV3	51201	102084	1.23%	0.078%
27	12.792	1625	1633	1640	rBV3	15325	36796	0.44%	0.028%
28	13.492	1744	1752	1757	rVV3	51583	119302	1.43%	0.092%
29	13.545	1757	1761	1766	rVV6	29026	56646	0.68%	0.043%
30	13.592	1766	1769	1774	rVB2	38846	55963	0.67%	0.043%
31	13.734	1789	1793	1798	rBV5	45104	72051	0.87%	0.055%
32	13.857	1808	1814	1818	rBV	165325	271384	3.26%	0.208%
33	13.904	1818	1822	1832	rVB	236944	369804	4.44%	0.284%
34	14.016	1832	1841	1850	rBV	2128179	3195960	38.40%	2.452%
35	14.292	1882	1888	1905	rBV	2578202	4154325	49.92%	3.187%
36	14.410	1905	1908	1914	rVB7	24156	40901	0.49%	0.031%

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Integrator: RTE
 Smoothing : OFF Filtering: 5
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 Start Thrs: 0.2 Max Peaks: 100
 Stop Thrs : 0 Peak Location: TOP

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

37	14.481	1914	1920	1924	rBV5	55218	92672	1.11%	0.071%
38	14.545	1924	1931	1934	rBV3	103400	149325	1.79%	0.115%
39	14.598	1934	1940	1947	rVV	2197081	3295792	39.60%	2.528%
40	14.751	1961	1966	1970	rBV5	26387	50581	0.61%	0.039%
41	14.798	1970	1974	1975	rBV2	49624	63570	0.76%	0.049%
42	14.863	1976	1985	1998	rVV2	1476002	3438761	41.32%	2.638%
43	14.986	2002	2006	2009	rBV	63682	87377	1.05%	0.067%
44	15.151	2030	2034	2038	rVB5	56880	85362	1.03%	0.065%
45	15.475	2084	2089	2096	rVB	423374	610560	7.34%	0.468%
46	15.592	2102	2109	2118	rBV	3422931	5079089	61.03%	3.896%
47	15.728	2126	2132	2143	rBV	627946	928948	11.16%	0.713%
48	15.875	2153	2157	2162	rVB	133204	180976	2.17%	0.139%
49	16.075	2188	2191	2195	rVB5	44382	59135	0.71%	0.045%
50	16.239	2216	2219	2224	rVB	51361	71142	0.85%	0.055%
51	16.310	2226	2231	2239	rBV	104773	191261	2.30%	0.147%
52	16.463	2249	2257	2261	rBV5	75910	161799	1.94%	0.124%
53	16.522	2264	2267	2276	rVB3	47488	84107	1.01%	0.065%
54	16.745	2301	2305	2314	rBV3	122698	225321	2.71%	0.173%
55	16.845	2318	2322	2327	rVB2	154294	214249	2.57%	0.164%
56	16.998	2344	2348	2356	rBV	155707	259774	3.12%	0.199%
57	17.239	2385	2389	2392	rBV3	128497	184356	2.22%	0.141%
58	17.304	2397	2400	2402	rBV2	62424	69804	0.84%	0.054%
59	17.345	2402	2407	2412	rBV	2530284	3852153	46.29%	2.955%
60	17.445	2418	2424	2433	rVB	3268610	4915731	59.07%	3.771%
61	17.557	2440	2443	2445	rBV4	52620	71175	0.86%	0.055%
62	17.586	2446	2448	2453	rVB2	91283	107450	1.29%	0.082%
63	17.704	2464	2468	2471	rBV	124629	169943	2.04%	0.130%
64	17.874	2494	2497	2502	rVB2	66807	79896	0.96%	0.061%
65	17.933	2505	2507	2511	rVB3	60212	61618	0.74%	0.047%
66	18.127	2537	2540	2546	rVB3	64150	89325	1.07%	0.069%
67	18.186	2547	2550	2555	rBV3	75896	144719	1.74%	0.111%
68	18.245	2555	2560	2570	rBV	1183879	1725309	20.73%	1.324%
69	18.357	2576	2579	2585	rVB2	148672	231161	2.78%	0.177%
70	18.539	2607	2610	2614	rBV4	86607	146432	1.76%	0.112%
71	18.586	2614	2618	2624	rVB4	107802	179199	2.15%	0.137%
72	19.033	2690	2694	2698	rBV3	120435	153109	1.84%	0.117%
73	19.204	2719	2723	2728	rVB2	205644	255106	3.07%	0.196%
74	19.257	2728	2732	2736	rBV	330899	430776	5.18%	0.330%
75	19.374	2747	2752	2754	rBV3	124671	211980	2.55%	0.163%
76	19.404	2754	2757	2759	rVV2	199039	276582	3.32%	0.212%

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77	19.439	2759	2763	2771	rVB2	641931	940704	11.30%	0.722%
78	19.521	2774	2777	2785	rVV	163480	282430	3.39%	0.217%
79	19.598	2787	2790	2797	rVB3	163387	233808	2.81%	0.179%
80	19.739	2808	2814	2820	rBV	4170583	5698900	68.48%	4.372%
81	20.251	2897	2901	2903	rBV3	353083	494562	5.94%	0.379%
82	20.286	2903	2907	2911	rVB	587181	806808	9.69%	0.619%
83	20.445	2930	2934	2942	rVB3	421934	734005	8.82%	0.563%
84	20.615	2958	2963	2967	rBV2	3530083	5850025	70.29%	4.488%
85	20.657	2967	2970	2977	rVB2	1233149	2065314	24.82%	1.584%
86	20.880	3006	3008	3012	rVB3	242152	272307	3.27%	0.209%
87	20.980	3021	3025	3030	rVB	1287659	1561584	18.76%	1.198%
88	21.115	3044	3048	3053	rBV	2566276	3374924	40.55%	2.589%
89	21.239	3065	3069	3073	rBV	2129019	2553332	30.68%	1.959%
90	21.339	3082	3086	3088	rBV3	639480	948114	11.39%	0.727%
91	21.533	3114	3119	3128	rBV	2939627	3975301	47.77%	3.050%
92	22.109	3214	3217	3221	rBV	634372	773608	9.30%	0.593%
93	22.174	3223	3228	3239	rVB	2643382	4049888	48.66%	3.107%
94	22.662	3307	3311	3326	rVB	984282	1794034	21.56%	1.376%
95	23.227	3402	3407	3411	rBV	2213025	3528603	42.40%	2.707%
96	23.274	3411	3415	3428	rVB	2654494	4584937	55.09%	3.517%
97	23.798	3498	3504	3510	rBV2	2729333	5389768	64.76%	4.135%
98	23.956	3525	3531	3542	rVB	2036003	4323714	51.95%	3.317%
99	24.603	3635	3641	3650	rVB	3925626	8322537	100.00%	6.384%
100	24.698	3652	3657	3670	rVB	1495179	3340935	40.14%	2.563%

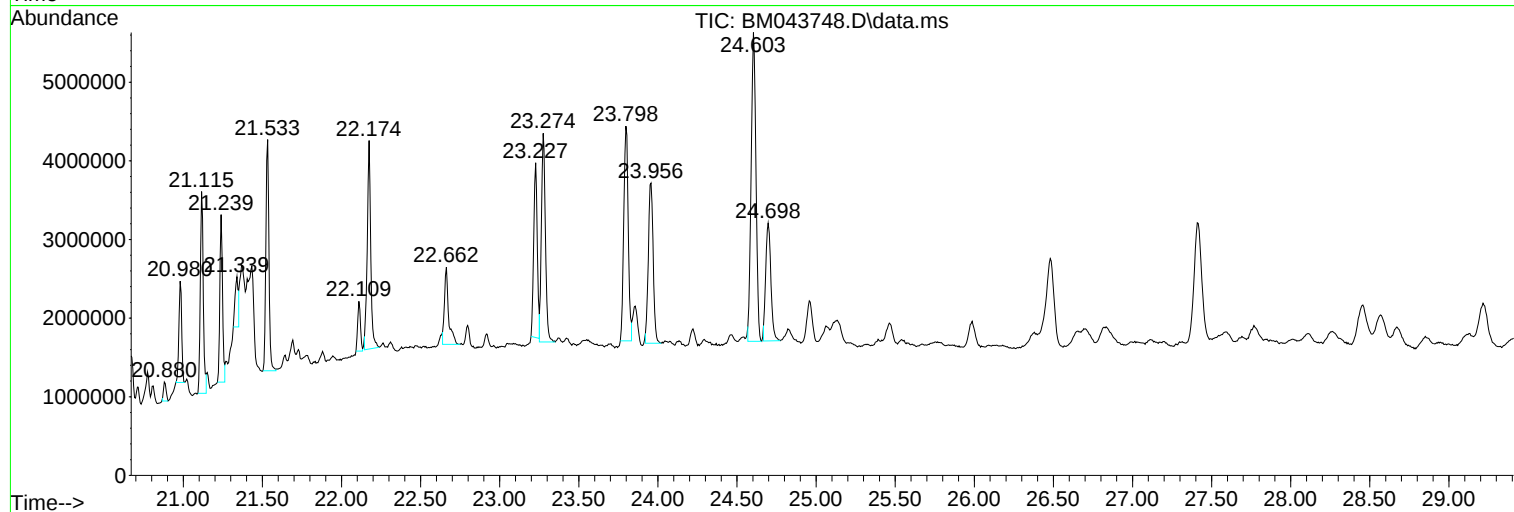
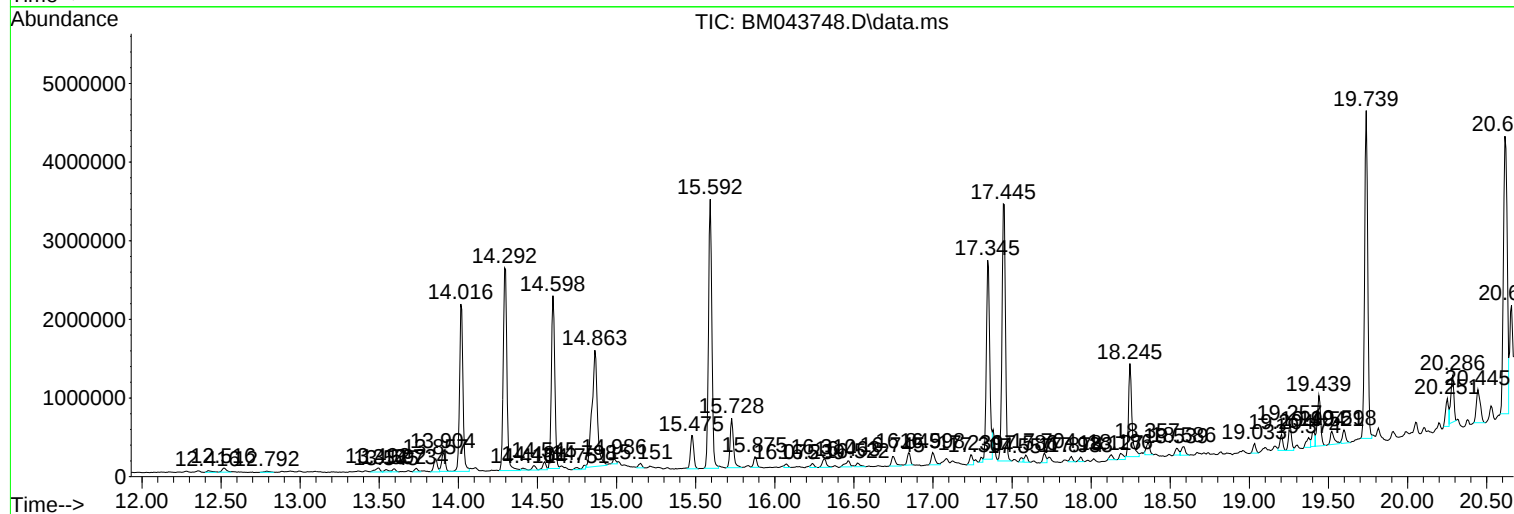
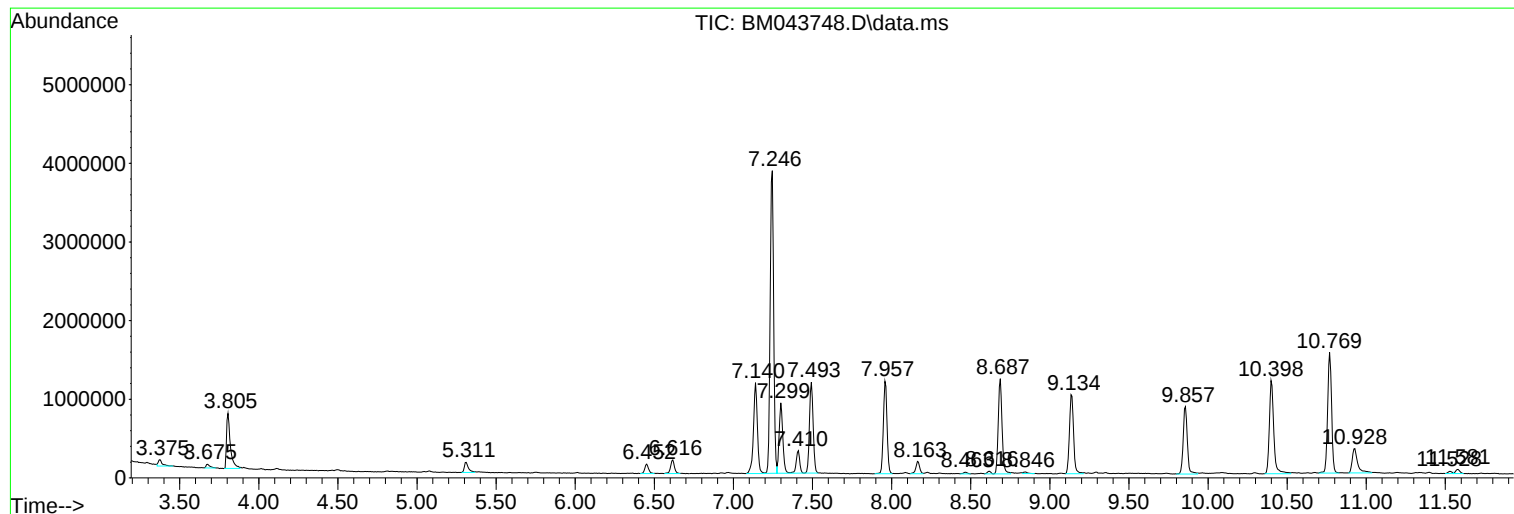
Sum of corrected areas: 130357780

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

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



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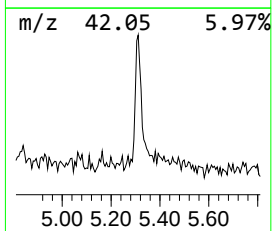
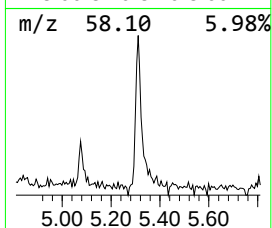
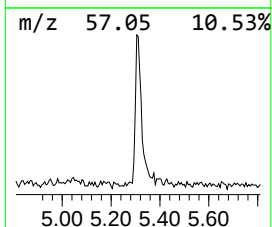
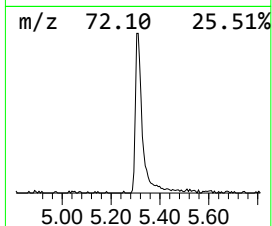
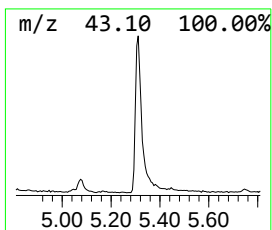
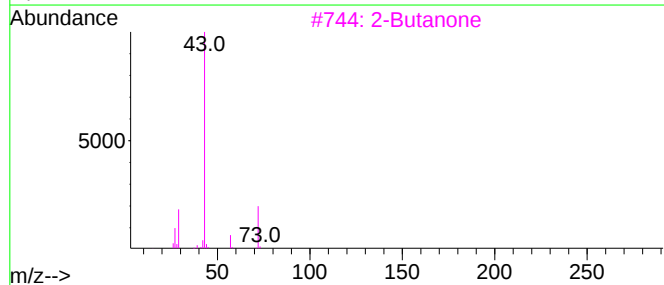
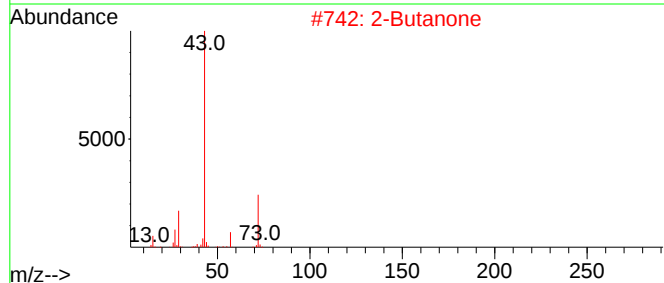
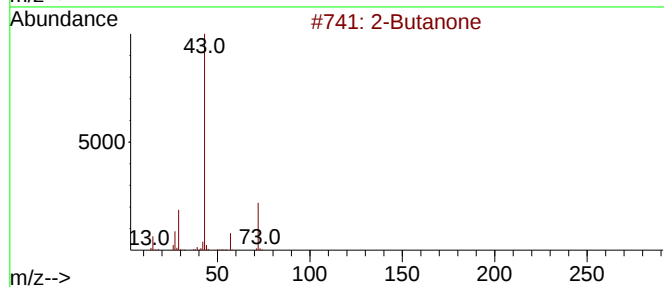
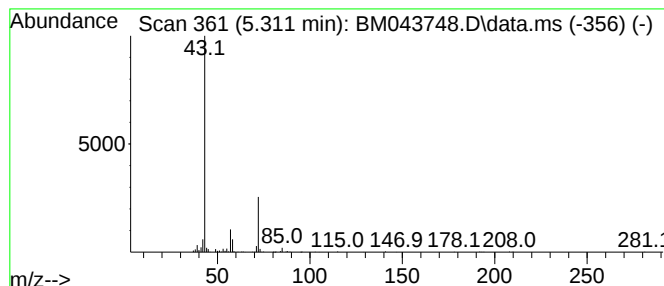
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM121523.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 2-Butanone Concentration Rank 24

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.311	2.49 ng/ul	236870	1,4-Dichlorobenzene-d4	7.957

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Butanone			72	C4H8O	000078-93-3	64
2	2-Butanone			72	C4H8O	000078-93-3	64
3	2-Butanone			72	C4H8O	000078-93-3	64
4	2-Butanone			72	C4H8O	000078-93-3	64
5	Propanal, 2-methyl-			72	C4H8O	000078-84-2	50



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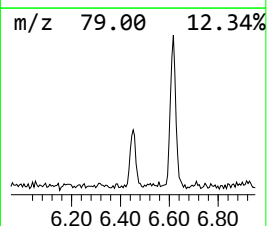
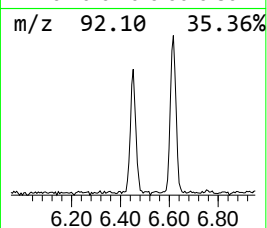
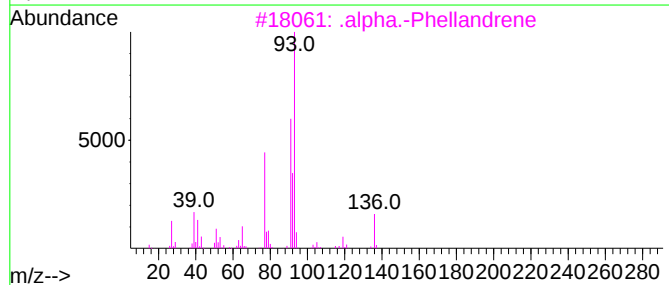
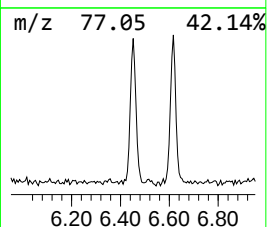
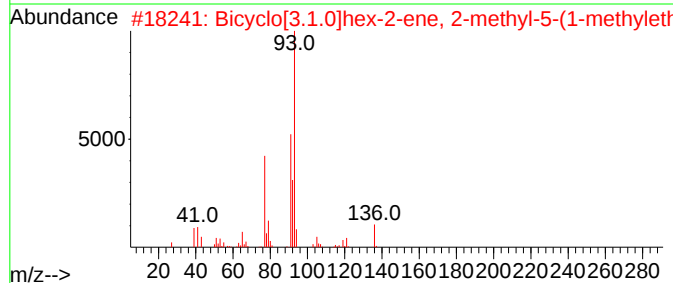
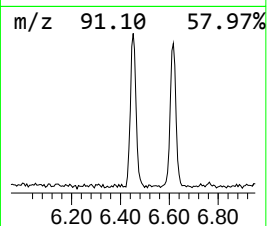
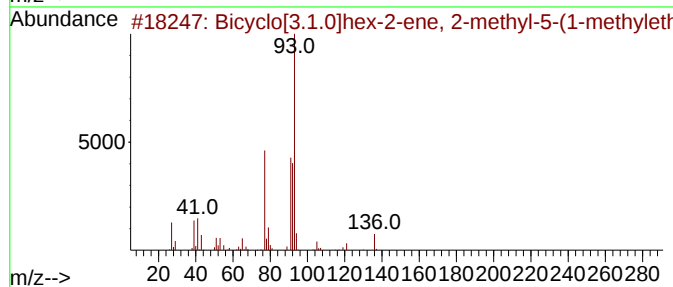
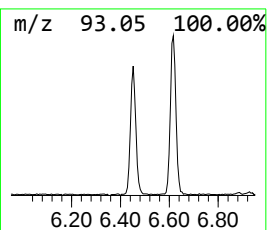
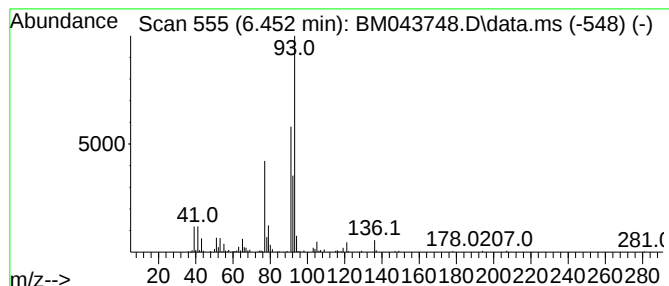
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM121523.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 Bicyclo[3.1.0]hex-2-ene, 2-... Concentration Rank 27

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.452	2.04 ng/ul	194464	1,4-Dichlorobenzene-d4	7.957

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Bicyclo[3.1.0]hex-2-ene, 2-methy...	136	C10H16	002867-05-2	91
2			Bicyclo[3.1.0]hex-2-ene, 2-methy...	136	C10H16	002867-05-2	91
3			.alpha.-Phellandrene	136	C10H16	000099-83-2	80
4			.alpha.-Phellandrene	136	C10H16	000099-83-2	78
5			Bicyclo[3.1.0]hex-2-ene, 2-methy...	136	C10H16	002867-05-2	78



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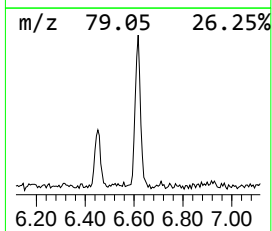
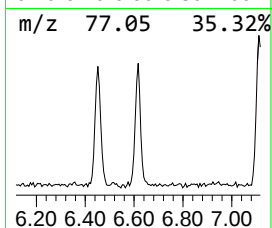
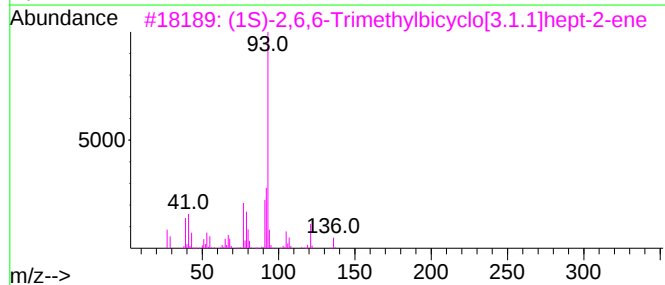
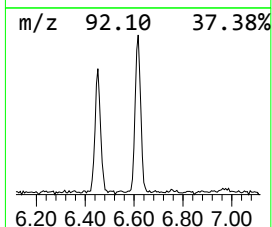
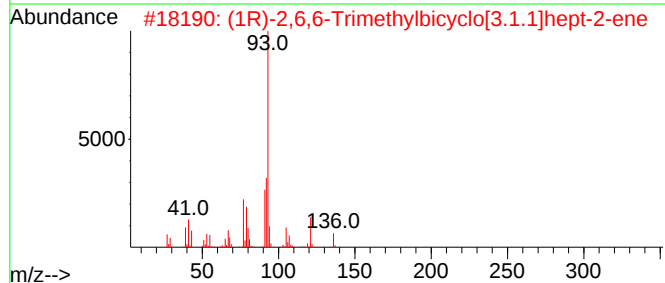
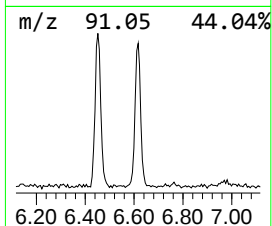
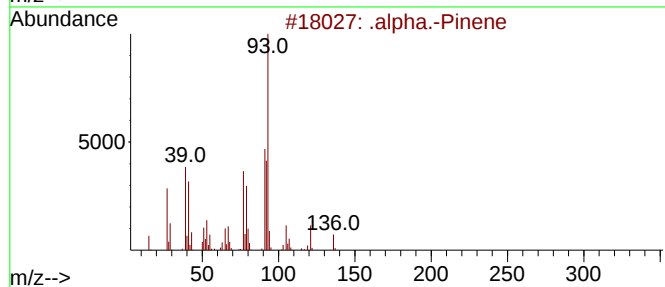
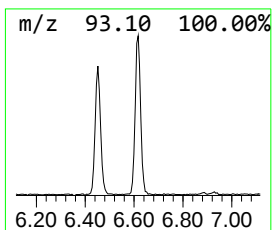
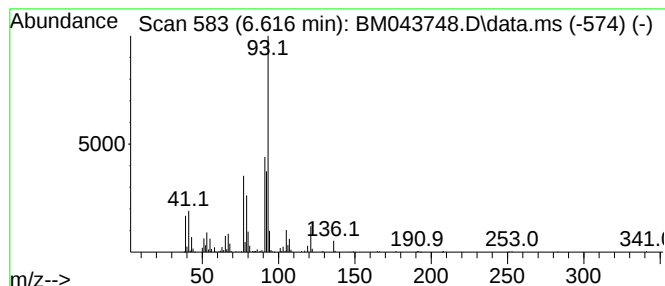
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Quant Title : SVOA CALIBRATION

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

Peak Number 3 .alpha.-Pinene Concentration Rank 21

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.616	3.11 ng/ul	295900	1,4-Dichlorobenzene-d4	7.957

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		.alpha.-Pinene	136	C10H16	000080-56-8	95
2		(1R)-2,6,6-Trimethylbicyclo[3.1.1]hept-2-ene	136	C10H16	007785-70-8	94
3		(1S)-2,6,6-Trimethylbicyclo[3.1.1]hept-2-ene	136	C10H16	007785-26-4	91
4		3-Carene	136	C10H16	013466-78-9	91
5		(+)-3-Carene	136	C10H16	000498-15-7	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM010224\
Data File : BM043748.D
Acq On : 04 Jan 2024 00:09
Operator : MA/JU
Sample : 05951-04
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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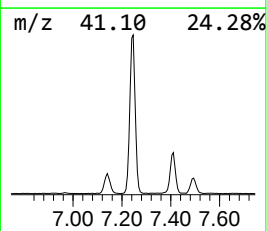
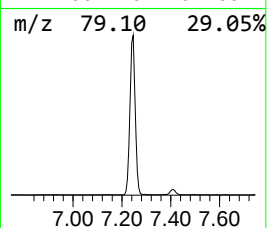
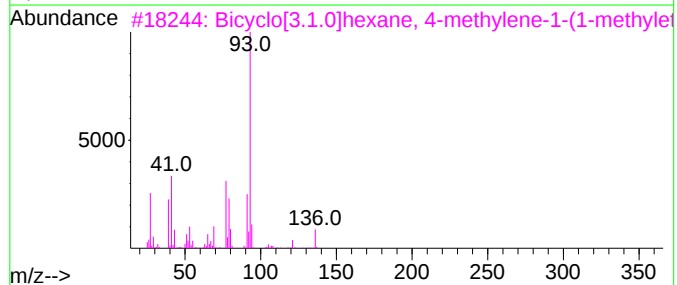
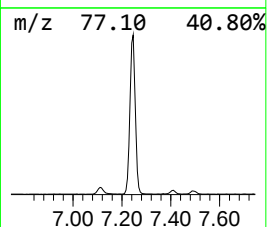
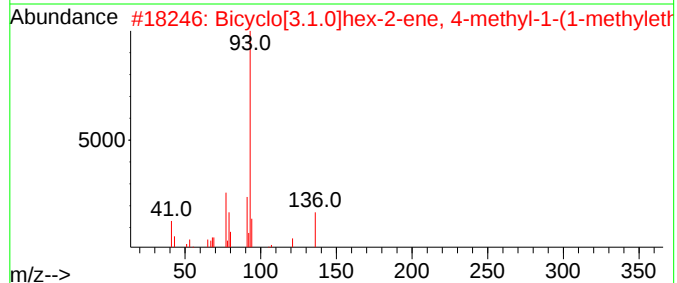
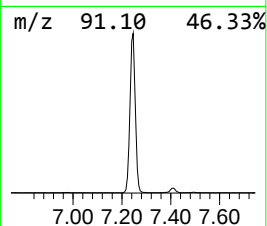
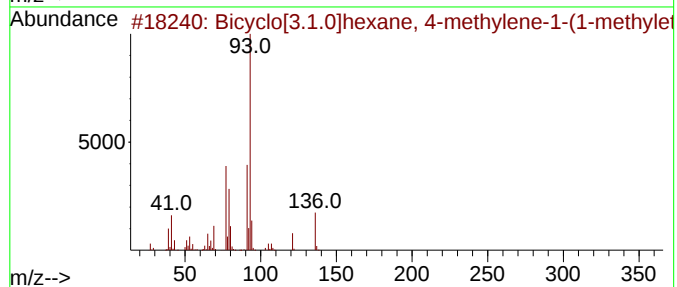
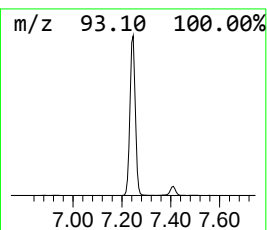
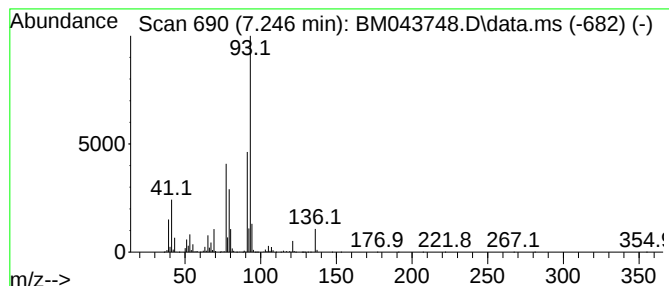
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 Bicyclo[3.1.0]hexane, 4-met... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.246	62.25 ng/ul	5927100	1,4-Dichlorobenzene-d4	7.957

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Bicyclo[3.1.0]hexane, 4-methylen...	136	C10H16	003387-41-5	97
2			Bicyclo[3.1.0]hex-2-ene, 4-methy...	136	C10H16	028634-89-1	91
3			Bicyclo[3.1.0]hexane, 4-methylen...	136	C10H16	003387-41-5	91
4			.beta.-Phellandrene	136	C10H16	000555-10-2	90
5			Bicyclo[3.1.0]hexane, 4-methylen...	136	C10H16	003387-41-5	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM010224\
Data File : BM043748.D
Acq On : 04 Jan 2024 00:09
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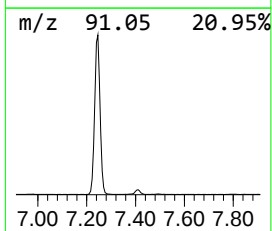
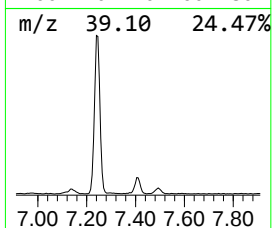
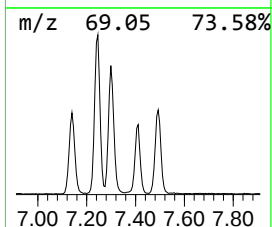
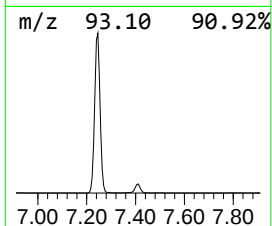
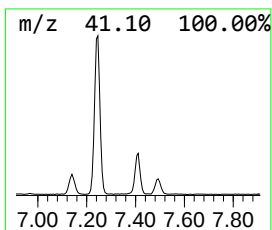
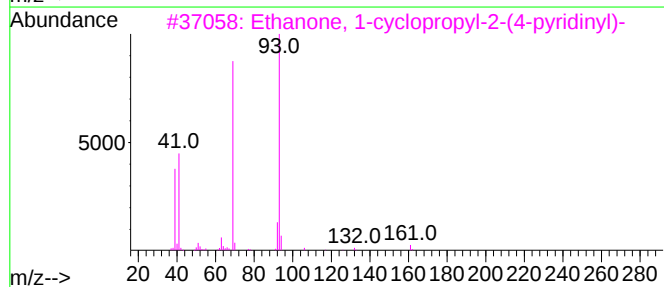
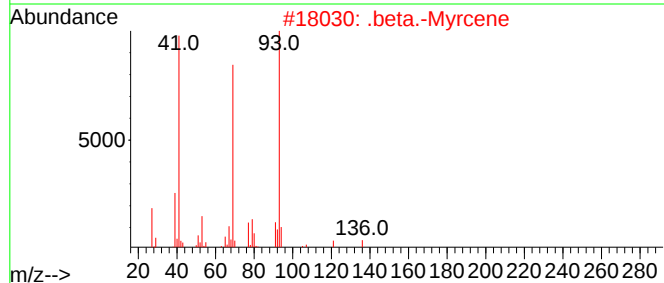
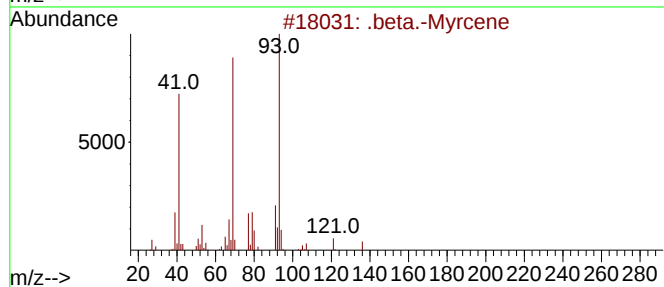
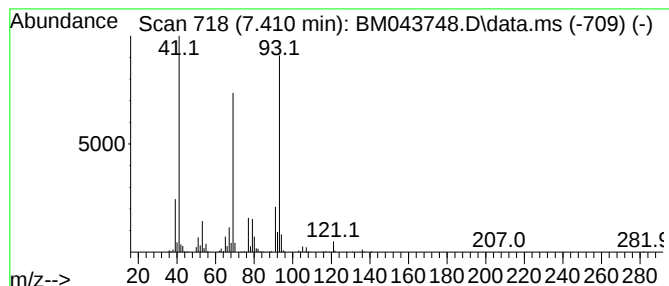
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 .beta.-Myrcene Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.410	4.59 ng/ul	436557	1,4-Dichlorobenzene-d4	7.957

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	.	beta.-Myrcene	136	C10H16	000123-35-3	90	
2	.	beta.-Myrcene	136	C10H16	000123-35-3	83	
3	Ethanone, 1-cyclopropyl-2-(4-pyr...	161	C10H11NO	006580-95-6	59		
4	.	beta.-Pinene	136	C10H16	000127-91-3	52	
5	.	beta.-Myrcene	136	C10H16	000123-35-3	50	



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM010224\
Data File : BM043748.D
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Sample : 05951-04
Misc :
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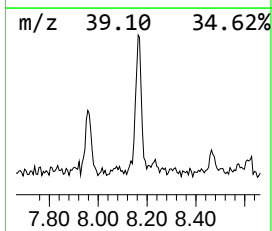
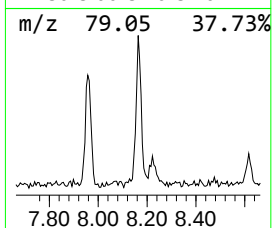
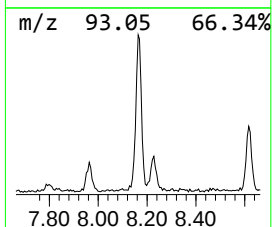
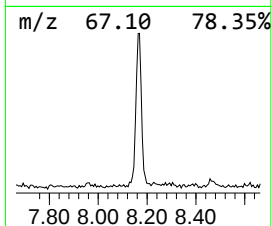
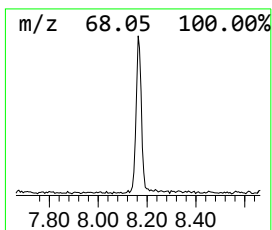
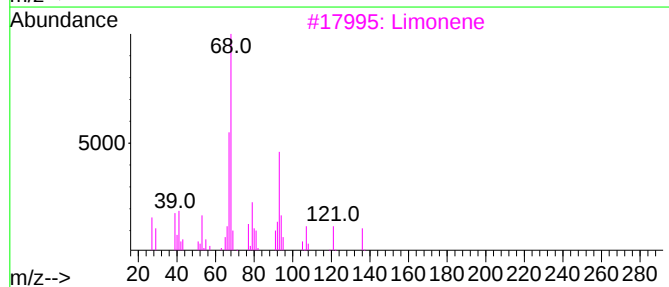
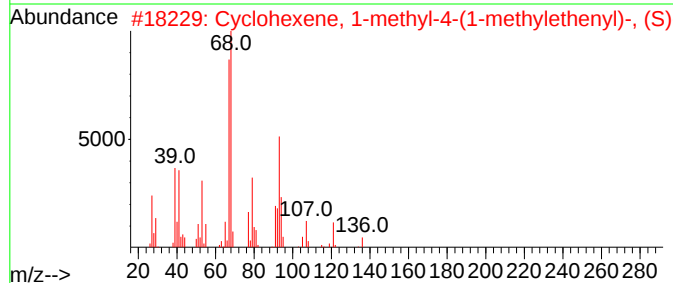
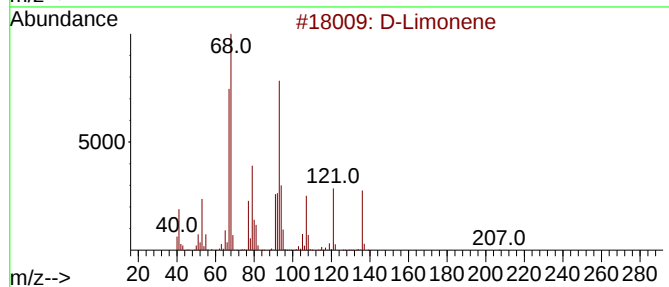
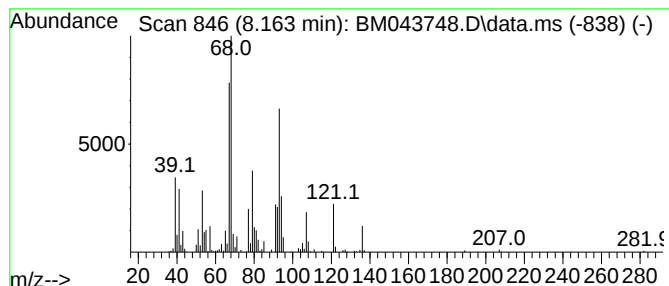
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 D-Limonene Concentration Rank 22

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.163	2.82 ng/ul	268728	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			Cyclohexene, 1-methyl-4-(1-methy...	136	C10H16	005989-54-8	94
3			Limonene	136	C10H16	000138-86-3	90
4			Cyclohexene, 1-methyl-5-(1-methy...	136	C10H16	013898-73-2	90
5			Limonene	136	C10H16	000138-86-3	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM010224\
Data File : BM043748.D
Acq On : 04 Jan 2024 00:09
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Sample : 05951-04
Misc :
ALS Vial : 16 Sample Multiplier: 1

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ClientSampleId :
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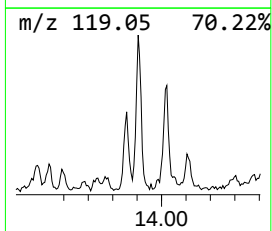
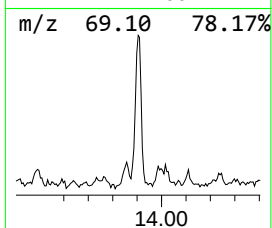
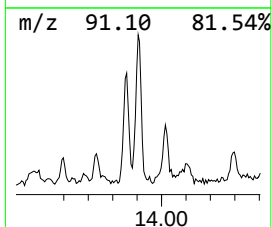
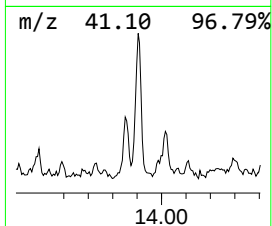
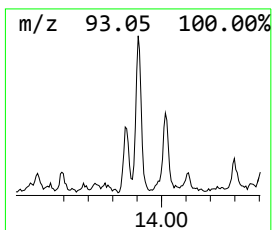
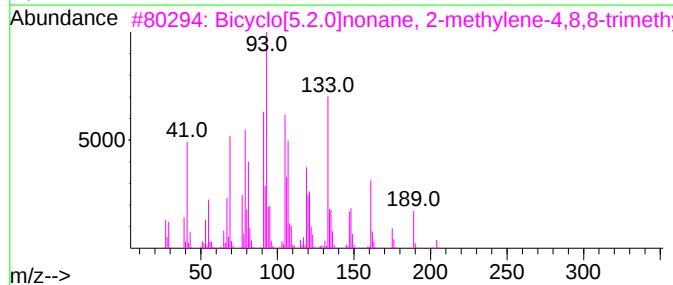
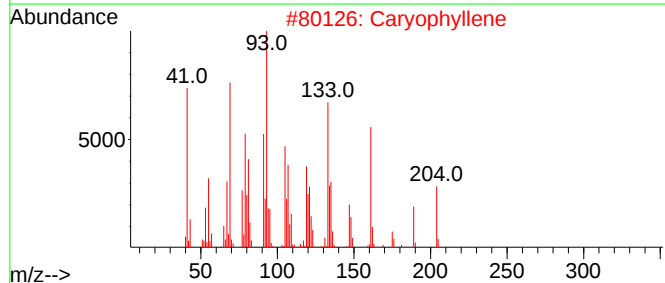
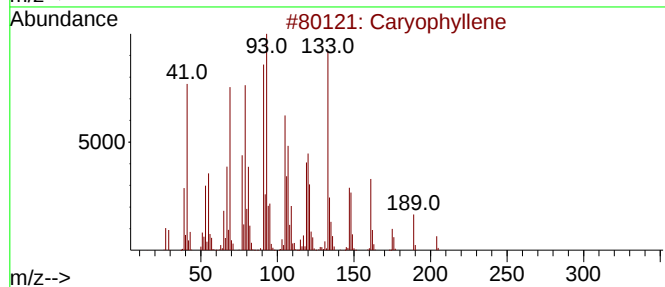
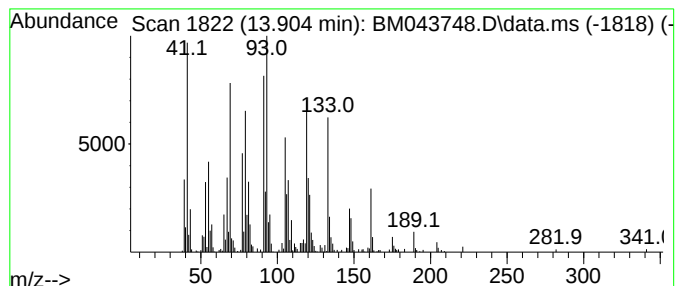
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 Caryophyllene Concentration Rank 25

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.904	2.24 ng/ul	369804	Acenaphthene-d10	14.598

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Caryophyllene	204	C15H24	000087-44-5	99
2			Caryophyllene	204	C15H24	000087-44-5	93
3			Bicyclo[5.2.0]nonane, 2-methylen...	204	C15H24	242794-76-9	91
4			Bicyclo[7.2.0]undec-4-ene, 4,11,...	204	C15H24	013877-93-5	91
5			Bicyclo[7.2.0]undec-4-ene, 4,11,...	204	C15H24	000118-65-0	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM010224\
Data File : BM043748.D
Acq On : 04 Jan 2024 00:09
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Sample : 05951-04
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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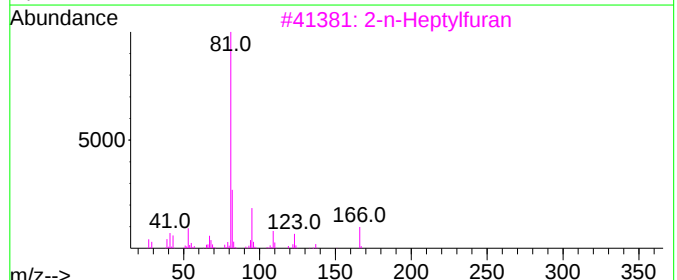
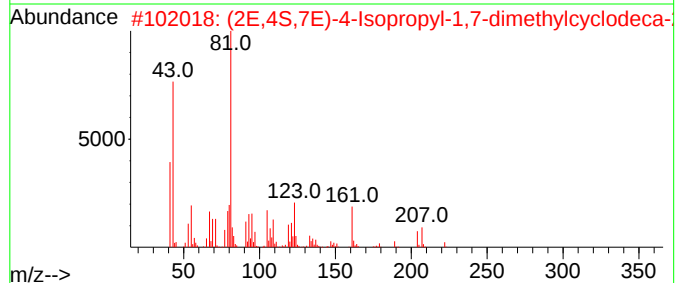
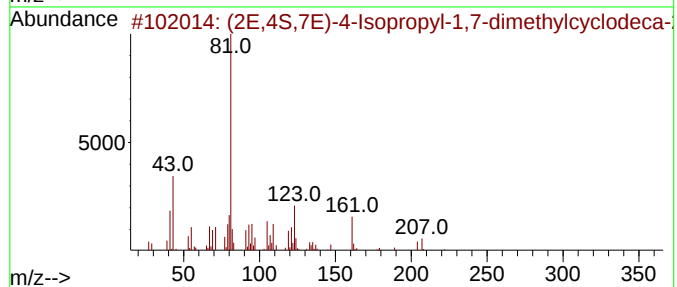
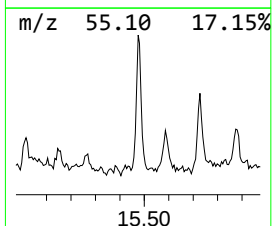
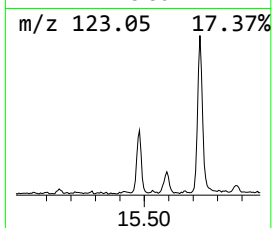
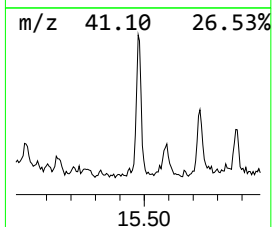
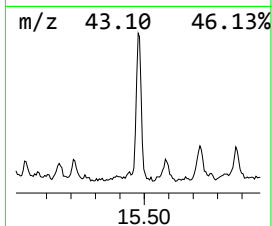
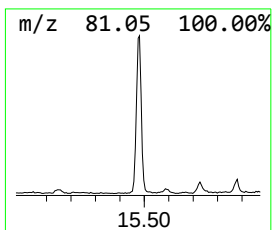
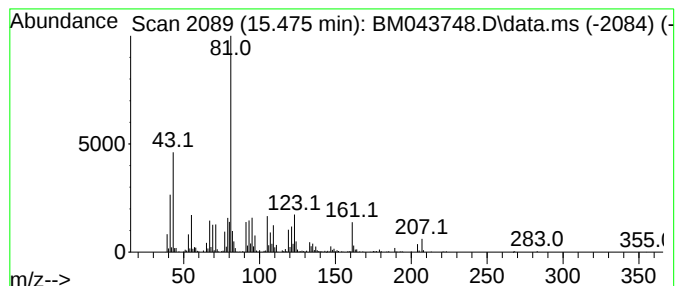
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 (2E,4S,7E)-4-Isopropyl-1,7-... Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.475	3.71 ng/ul	610560	Acenaphthene-d10	14.598

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			(2E,4S,7E)-4-Isopropyl-1,7-dimet...	222	C15H26O	198991-79-6	91
2			(2E,4S,7E)-4-Isopropyl-1,7-dimet...	222	C15H26O	198991-79-6	91
3			2-n-Heptylfuran	166	C11H18O	003777-71-7	50
4			16-Heptadecyn-4-one, 1,2-dihydroxy-	282	C17H30O3	1379770-10-1	50
5			Fenchone	152	C10H16O	001195-79-5	49



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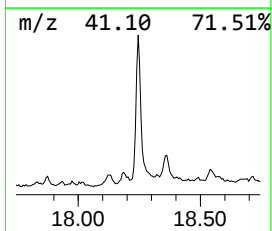
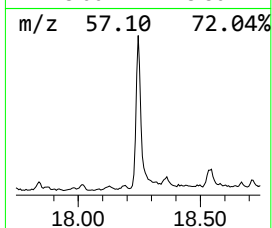
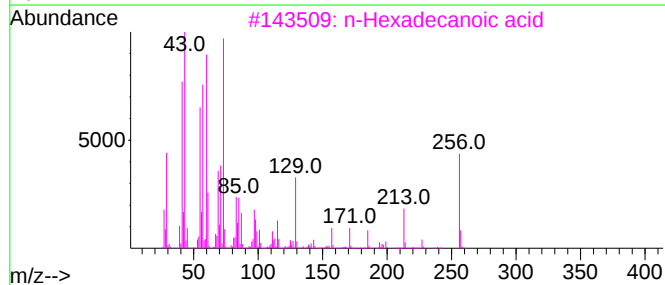
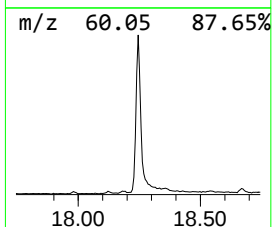
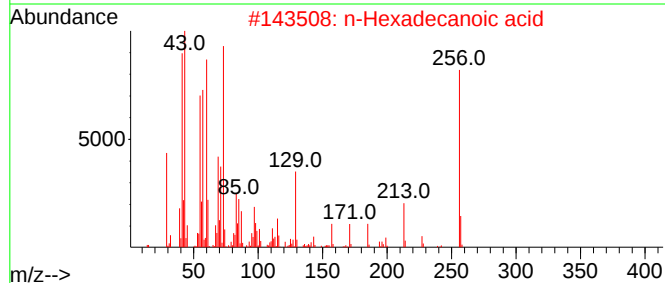
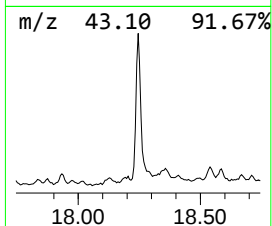
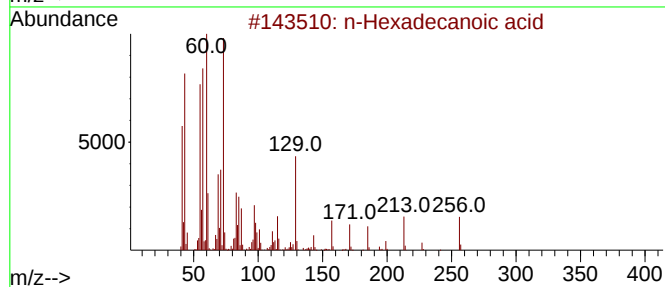
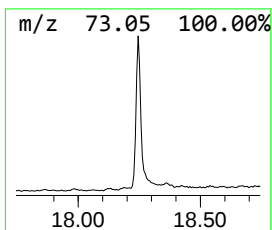
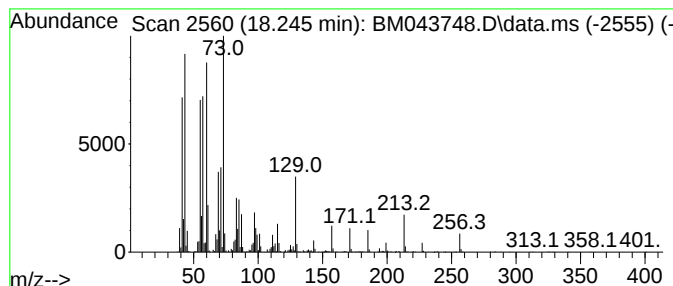
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 n-Hexadecanoic acid Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.	
18.245	8.96 ng/ul	1725310	Phenanthrene-d10	17.345	
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	98
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\
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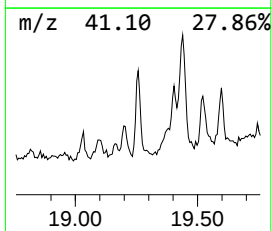
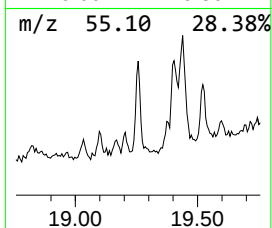
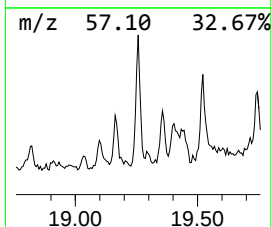
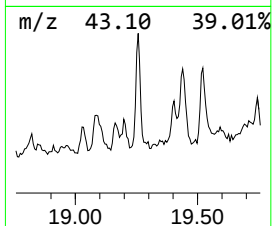
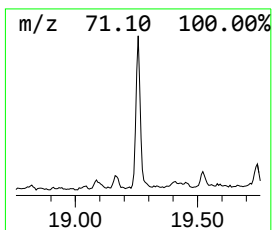
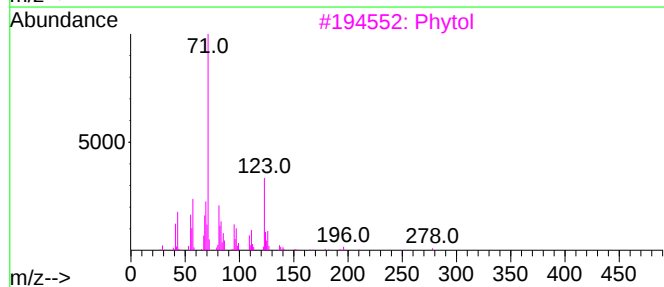
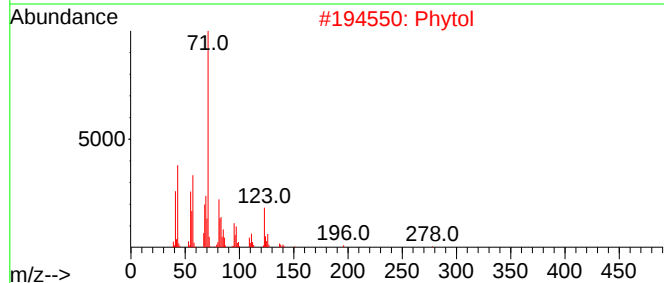
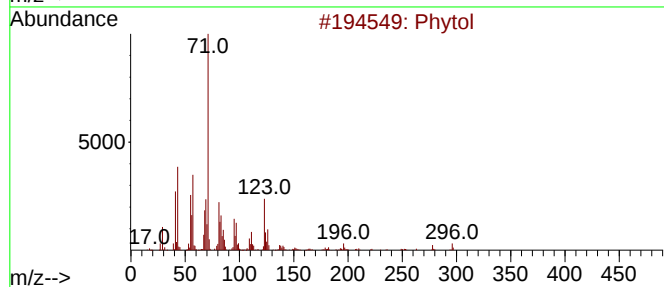
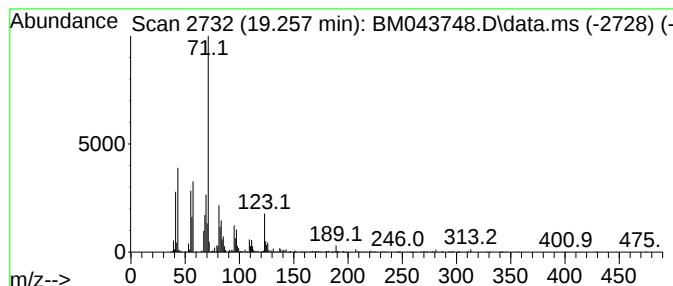
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 Phytol Concentration Rank 26

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.257	2.24 ng/ul	430776	Phenanthrene-d10	17.345

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phytol			296	C20H40O	000150-86-7	93
2	Phytol			296	C20H40O	000150-86-7	91
3	Phytol			296	C20H40O	000150-86-7	91
4	Phytol			296	C20H40O	000150-86-7	83
5	Butyric acid, 3-pentadecyl ester			298	C19H38O2	1000280-57-3	53



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM010224\
Data File : BM043748.D
Acq On : 04 Jan 2024 00:09
Operator : MA/JU
Sample : 05951-04
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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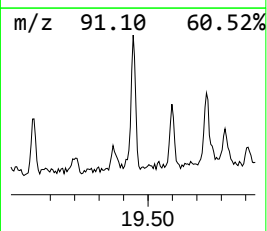
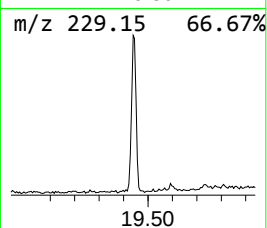
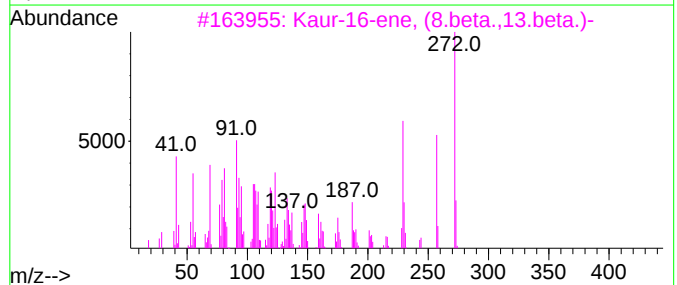
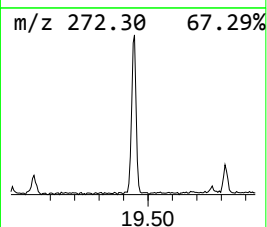
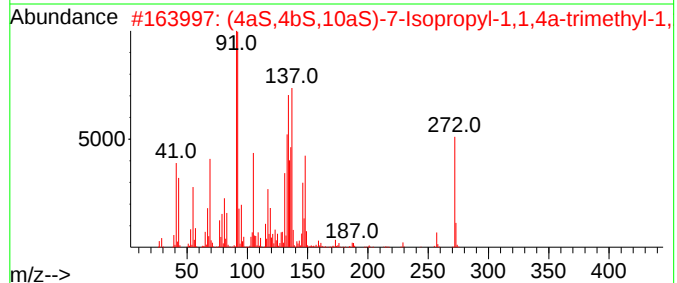
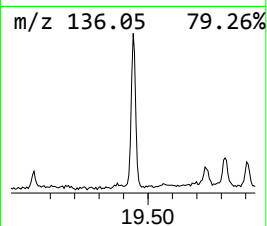
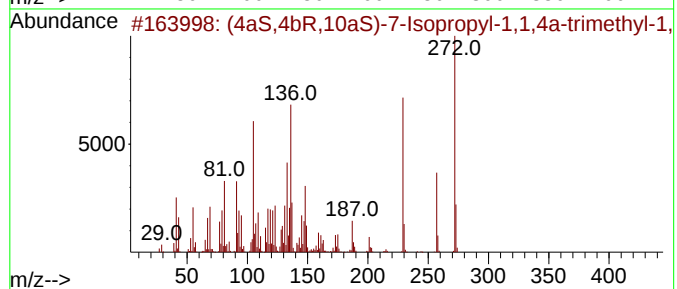
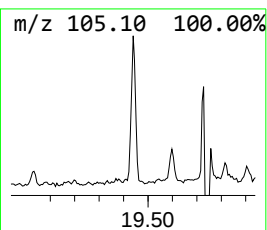
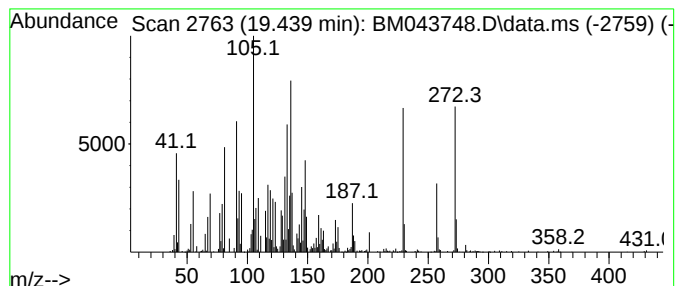
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Quant Title : SVOA CALIBRATION

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

Peak Number 11 (4aS,4bR,10aS)-7-Isopropyl-... Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.439	4.88 ng/ul	940704	Phenanthrene-d10	17.345

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			(4aS,4bR,10aS)-7-Isopropyl-1,1,4...	272	C20H32	035241-40-8	99
2			(4aS,4bS,10aS)-7-Isopropyl-1,1,4...	272	C20H32	122712-77-0	78
3			Kaur-16-ene, (8.beta.,13.beta.)-	272	C20H32	020070-61-5	55
4			Androst-5-en-4-one	272	C19H28O	013583-72-7	41
5			Benzocyclododecene, 2,3-diethyl-...	272	C20H32	061141-65-9	30



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM010224\
Data File : BM043748.D
Acq On : 04 Jan 2024 00:09
Operator : MA/JU
Sample : 05951-04
Misc :
ALS Vial : 16 Sample Multiplier: 1

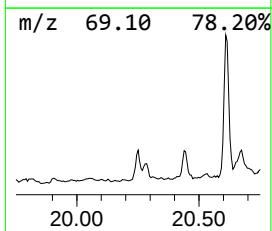
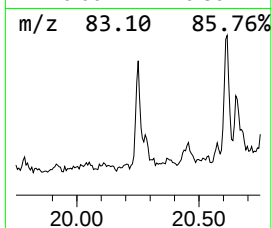
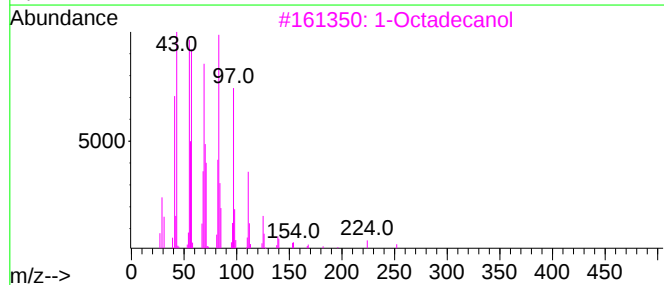
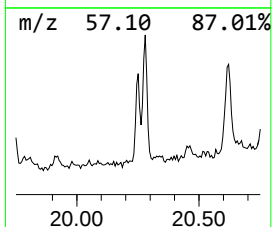
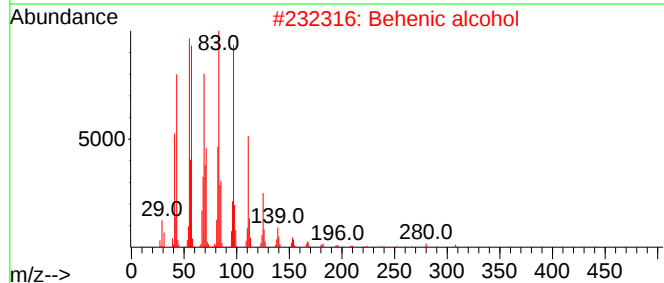
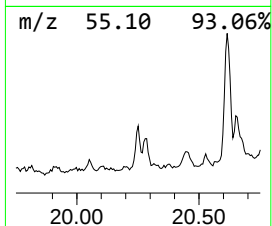
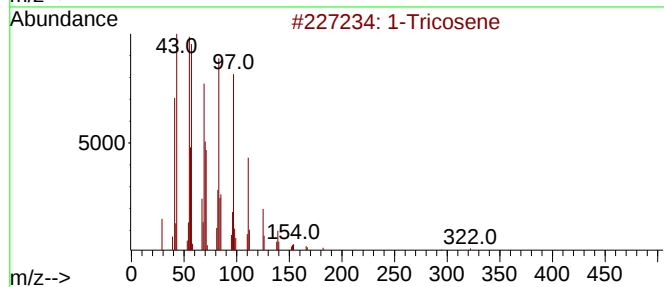
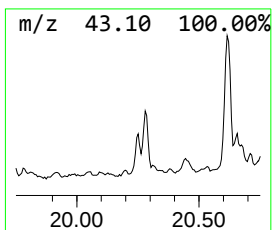
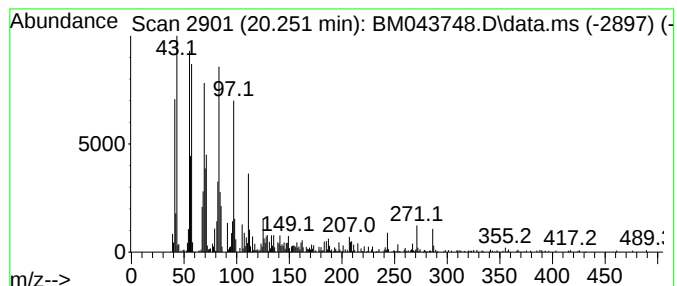
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ClientSampleId :
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Quant Title : SVOA CALIBRATION

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

Peak Number 12 (DEL) Alkane: Straight-Chai... Concentration Rank 23

R.T.	EstConc	Area	Relative to ISTD	R.T.	
20.251	2.49 ng/ul	494562	Chrysene-d12	21.533	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Tricosene	322	C23H46	018835-32-0	94
2	Behenic alcohol	326	C22H46O	000661-19-8	94
3	1-Octadecanol	270	C18H38O	000112-92-5	94
4	cis-1-Chloro-9-octadecene	286	C18H35Cl	016507-61-2	94
5	n-Nonadecanol-1	284	C19H40O	001454-84-8	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM010224\
Data File : BM043748.D
Acq On : 04 Jan 2024 00:09
Operator : MA/JU
Sample : 05951-04
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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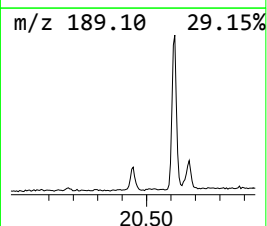
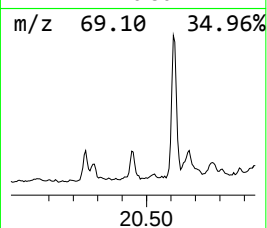
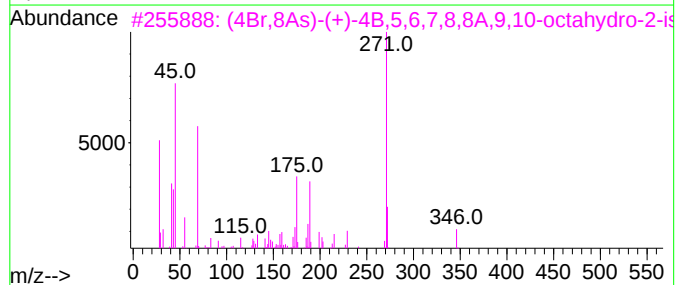
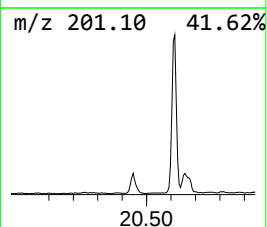
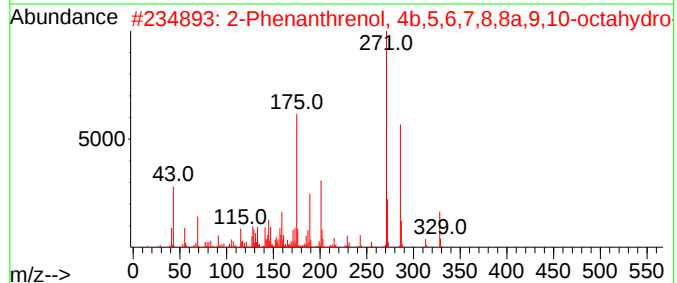
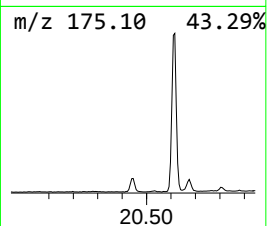
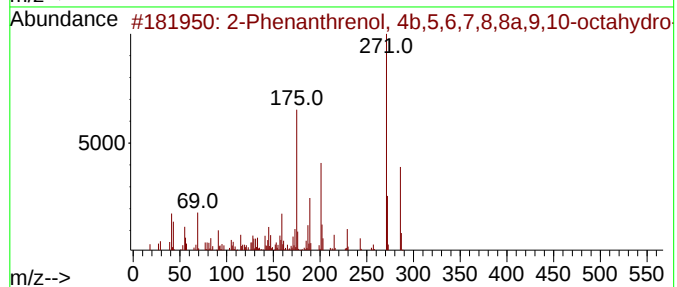
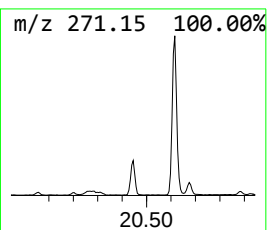
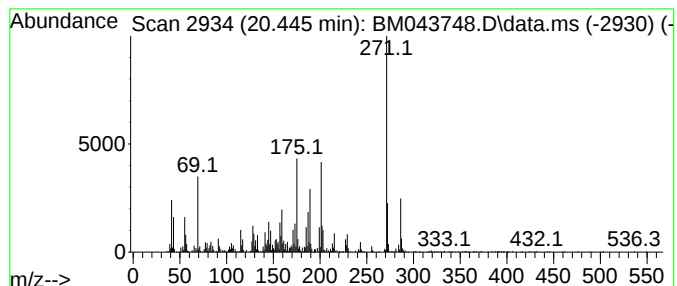
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Quant Title : SVOA CALIBRATION

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

Peak Number 14 2-Phenanthrenol, 4b,5,6,7,8... Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.445	3.69 ng/ul	734005	Chrysene-d12	21.533

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Phenanthrenol, 4b,5,6,7,8,8a,9...	286	C20H30O	000511-15-9	91
2			2-Phenanthrenol, 4b,5,6,7,8,8a,9...	328	C22H32O2	015340-82-6	58
3			(4Br,8As)-(+)-4B,5,6,7,8,8A,9,10...	346	C22H34O3	1000426-72-2	46
4			Silane, methylvinyl(2,4,4-trimet...	400	C21H44O3Si2	1000416-59-2	43
5			Podocarpa-8,11,13-trien-17-oic a...	344	C22H32O3	024067-43-4	43



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM010224\
Data File : BM043748.D
Acq On : 04 Jan 2024 00:09
Operator : MA/JU
Sample : 05951-04
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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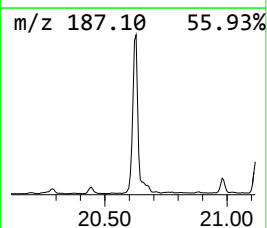
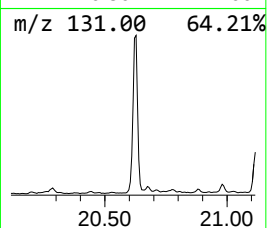
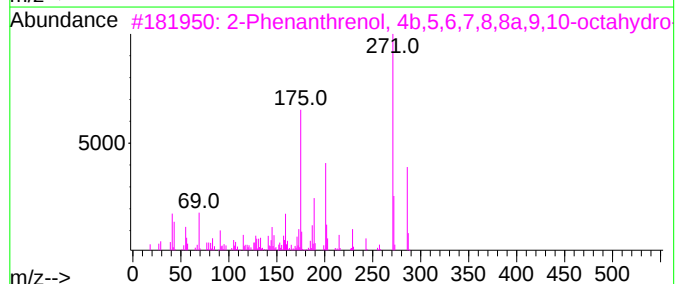
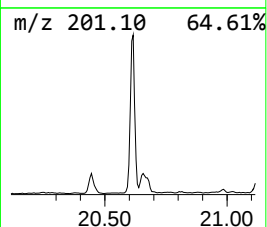
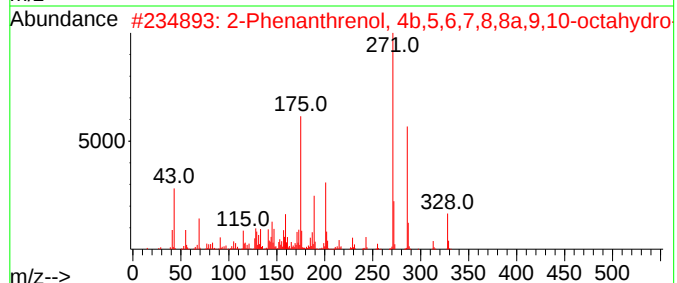
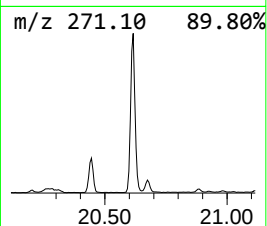
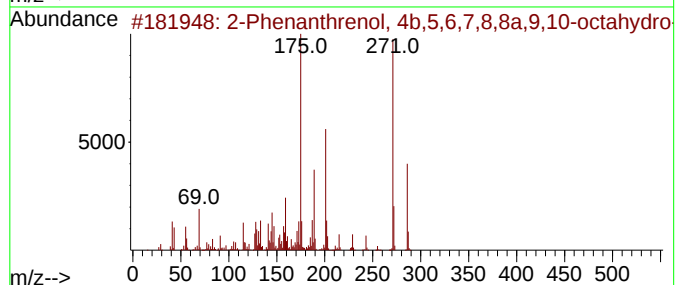
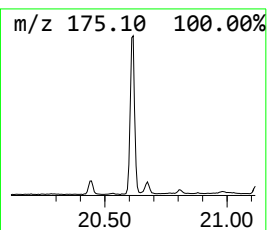
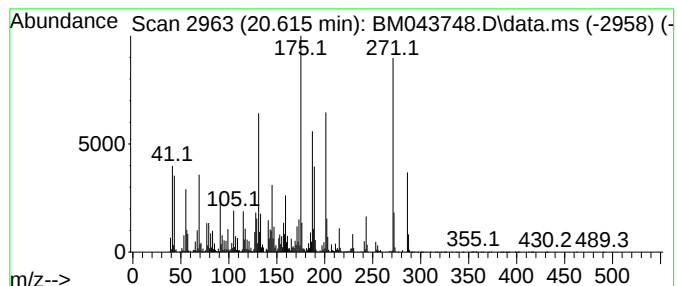
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Quant Title : SVOA CALIBRATION

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

Peak Number 15 Cycloisolongifolene, 9,10-d... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.616	29.43 ng/ul	5850030	Chrysene-d12	21.533

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Phenanthrenol, 4b,5,6,7,8,8a,9...	286	C20H300	000511-15-9	99		
2	2-Phenanthrenol, 4b,5,6,7,8,8a,9...	328	C22H3202	015340-82-6	68		
3	2-Phenanthrenol, 4b,5,6,7,8,8a,9...	286	C20H300	000511-15-9	68		
4	Cycloisolongifolene, 9,10-dehydro-	202	C15H22	1000156-81-6	53		
5	Pisiferol	302	C20H3002	024035-36-7	38		



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM010224\
Data File : BM043748.D
Acq On : 04 Jan 2024 00:09
Operator : MA/JU
Sample : 05951-04
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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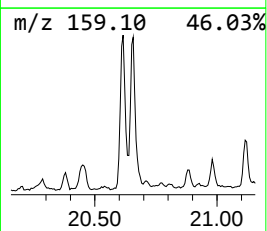
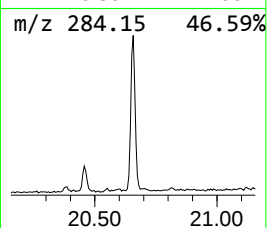
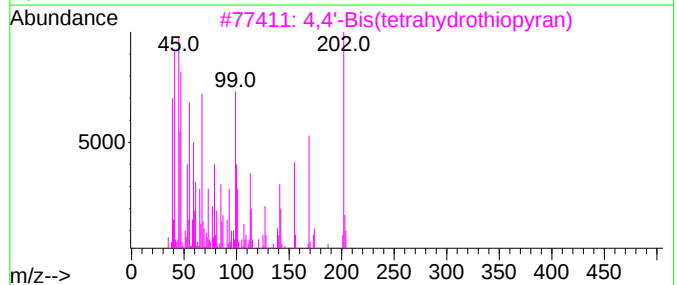
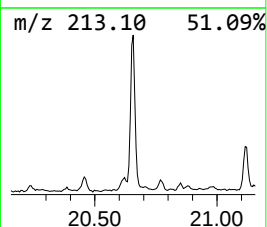
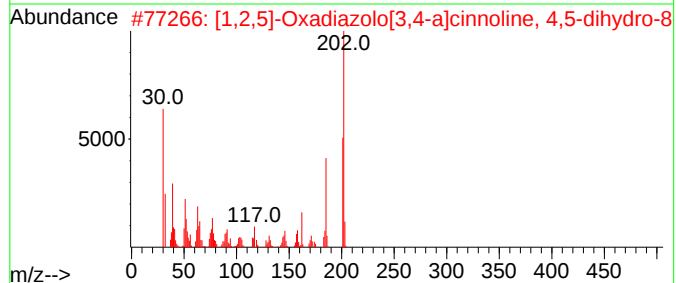
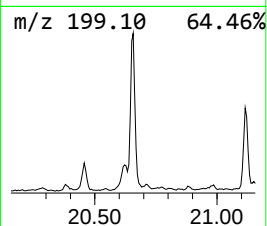
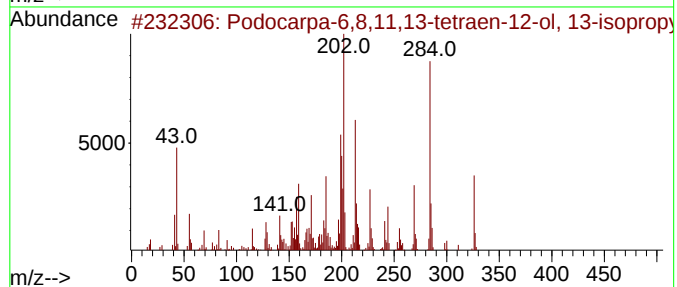
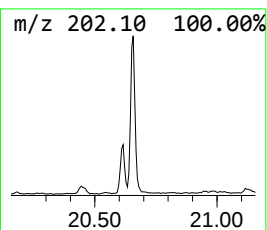
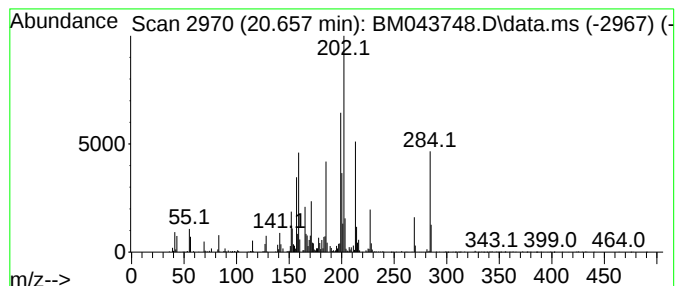
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Quant Title : SVOA CALIBRATION

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

Peak Number 16 unknown-01 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.657	10.39 ng/ul	2065310	Chrysene-d12	21.533

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Podocarpa-6,8,11,13-tetraen-12-o...	326	C22H30O2	022160-86-7	43
2			[1,2,5]-Oxadiazolo[3,4-a]cinnoli...	202	C10H10N4O	1000260-59-4	25
3			4,4'-Bis(tetrahydrothiopyran)	202	C10H18S2	116196-83-9	25
4			5-Hydroxy-3-methyl-1-phenylpyraz...	202	C11H10N2O2	060484-29-9	22
5			5-Hydroxy-3-methyl-1-phenylpyraz...	202	C11H10N2O2	060484-29-9	22



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM010224\
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Acq On : 04 Jan 2024 00:09
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Sample : 05951-04
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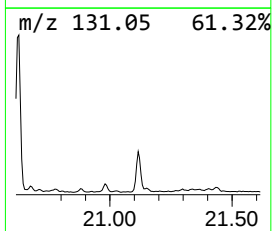
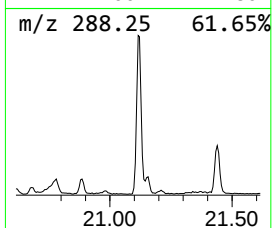
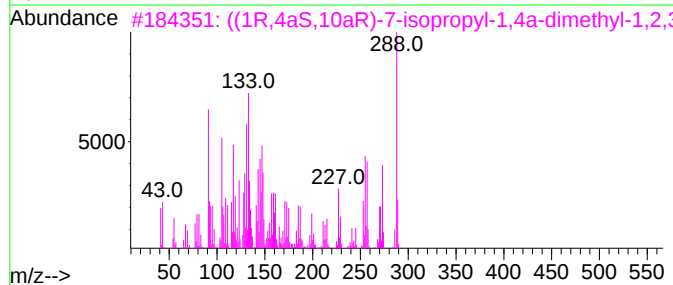
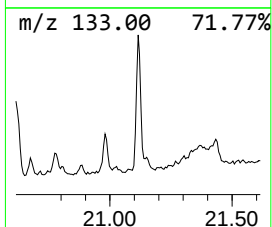
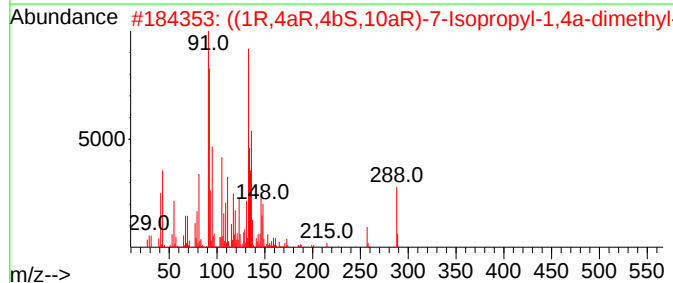
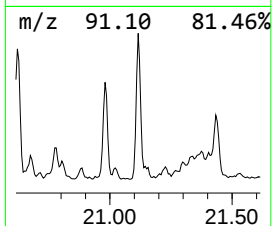
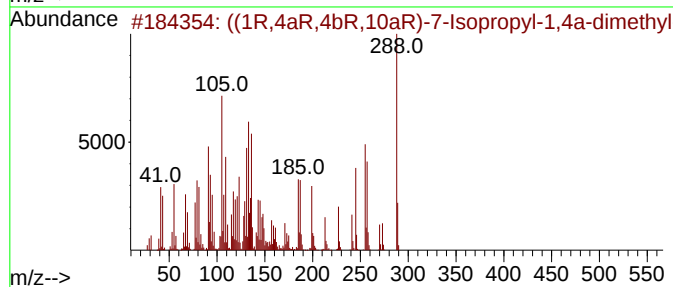
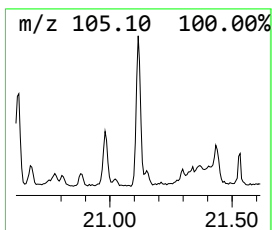
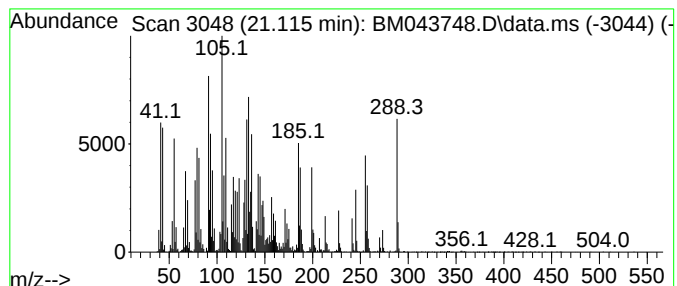
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ClientSampleId :
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Quant Title : SVOA CALIBRATION

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

Peak Number 18 ((1R,4aR,4bR,10aR)-7-Isopro... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.	
21.115	16.98 ng/ul	3374920	Chrysene-d12	21.533	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	((1R,4aR,4bR,10aR)-7-Isopropyl-1...	288	C20H32O	000666-84-2	99
2	((1R,4aR,4bS,10aR)-7-Isopropyl-1...	288	C20H32O	097640-46-5	38
3	((1R,4aS,10aR)-7-isopropyl-1,4a-...	288	C20H32O	1000439-86-6	35
4	3-Hydroxy-10,13-dimethyl-1,2,3,6...	288	C19H28O2	1000496-48-2	15
5	4'-Bromo-o-acetotoluidide	227	C9H10BrNO	024106-05-6	11



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM010224\
Data File : BM043748.D
Acq On : 04 Jan 2024 00:09
Operator : MA/JU
Sample : 05951-04
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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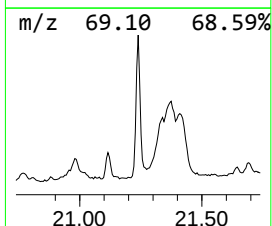
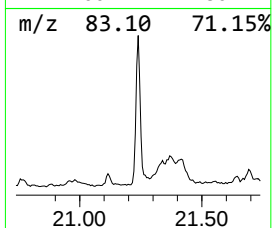
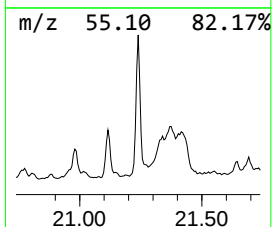
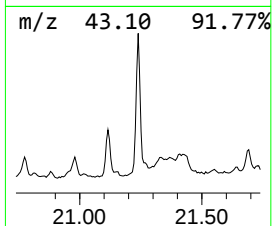
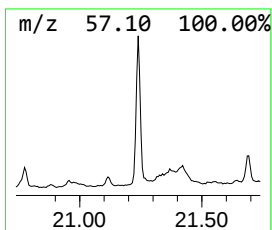
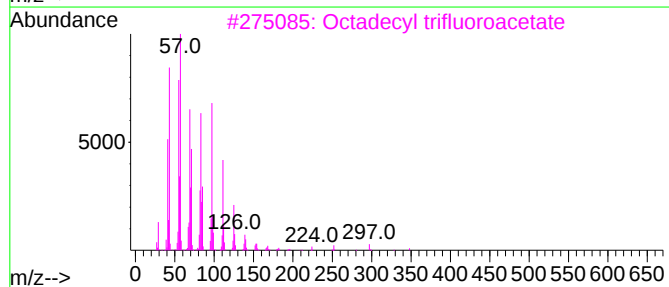
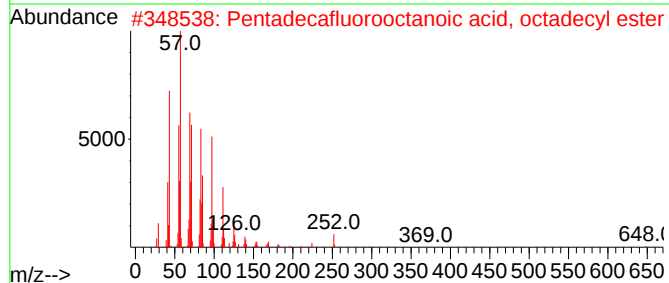
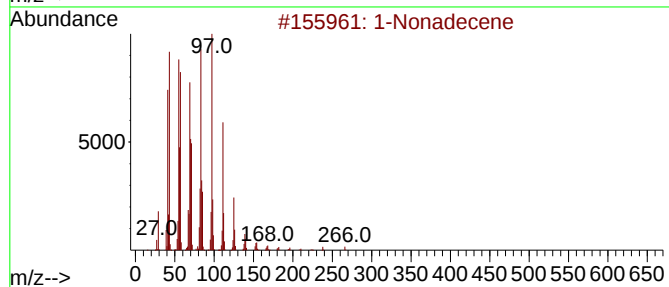
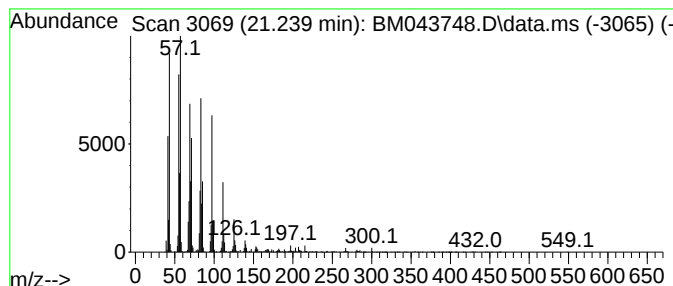
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Quant Title : SVOA CALIBRATION

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

Peak Number 19 (DEL) Alkane: Straight-Chai... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.239	12.85 ng/ul	2553330	Chrysene-d12	21.533

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Nonadecene			266	C19H38	018435-45-5	96
2	Pentadecafluorooctanoic acid, oc...			666	C26H37F15O2	1000406-04-8	95
3	Octadecyl trifluoroacetate			366	C20H37F3O2	079392-43-1	94
4	1-Hexacosene			364	C26H52	018835-33-1	94
5	Cyclotetracosane			336	C24H48	000297-03-0	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM010224\
Data File : BM043748.D
Acq On : 04 Jan 2024 00:09
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Sample : 05951-04
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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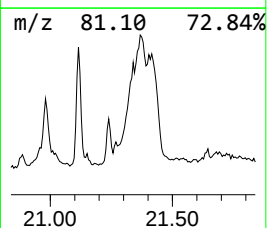
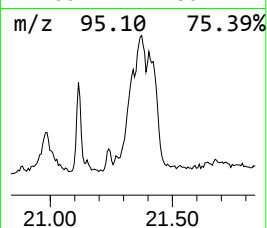
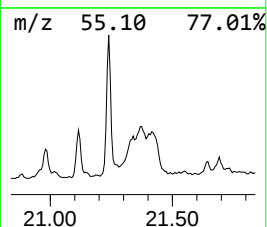
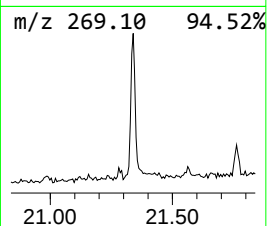
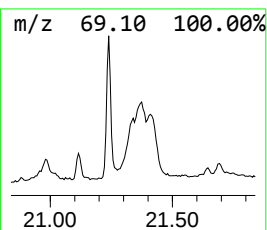
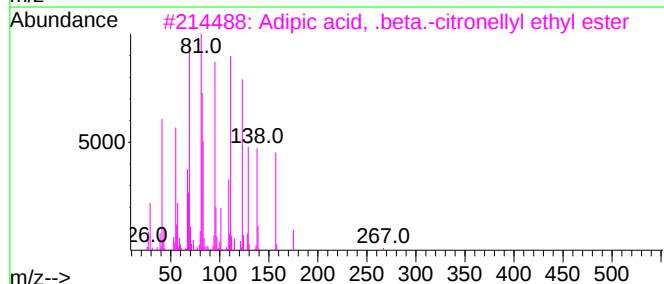
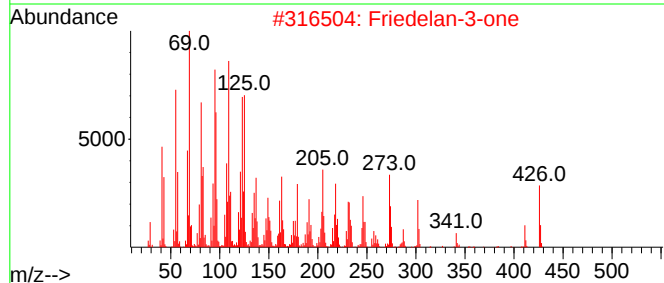
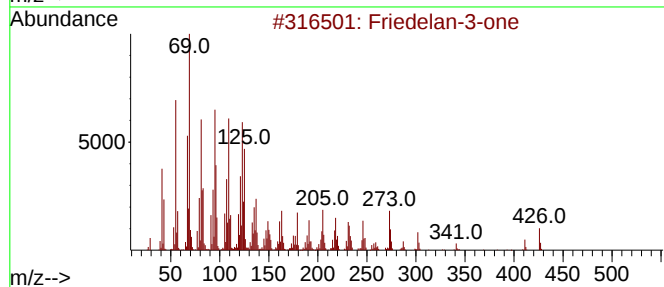
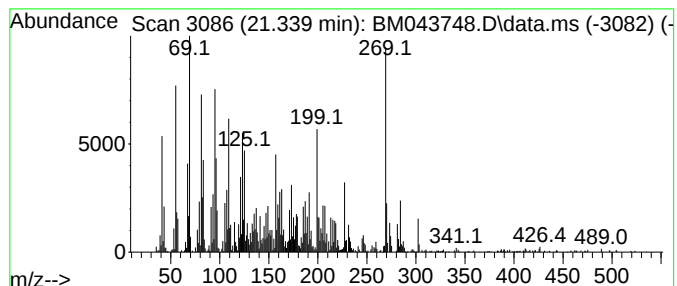
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Quant Title : SVOA CALIBRATION

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

Peak Number 20 unknown-02 Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.339	4.77 ng/ul	948114	Chrysene-d12	21.533

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Friedelan-3-one	426	C30H50O	000559-74-0	89
2			Friedelan-3-one	426	C30H50O	000559-74-0	53
3			Adipic acid, .beta.-citronellyl ...	312	C18H32O4	1000353-76-4	27
4			Phenol, 4,4'-(1-methylethylidene...	284	C19H24O2	005613-46-7	25
5			Friedelan-3-one	426	C30H50O	000559-74-0	25



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM010224\
Data File : BM043748.D
Acq On : 04 Jan 2024 00:09
Operator : MA/JU
Sample : 05951-04
Misc :
ALS Vial : 16 Sample Multiplier: 1

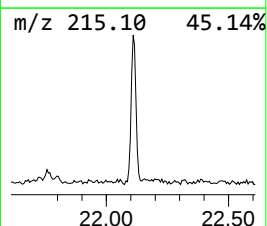
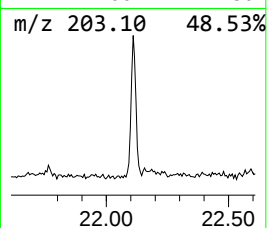
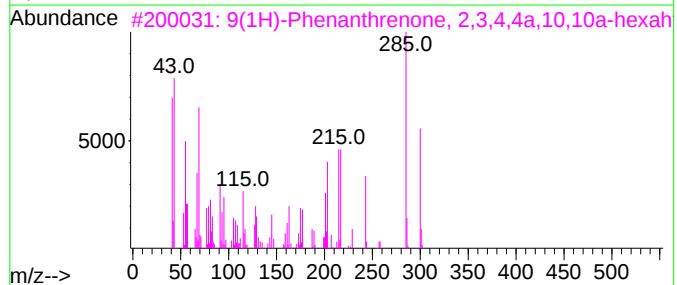
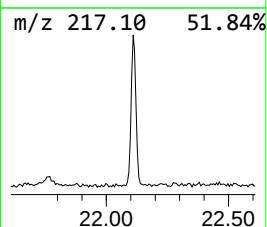
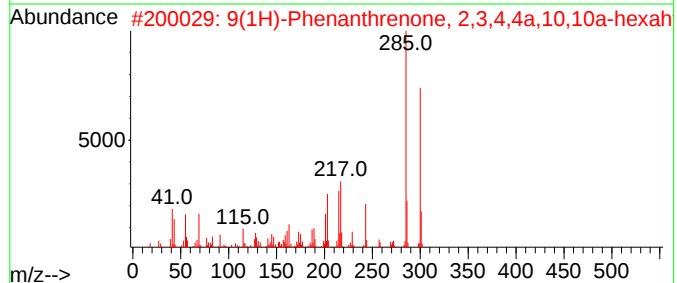
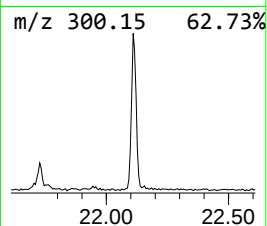
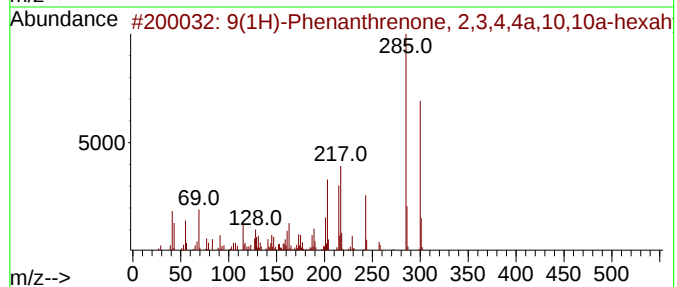
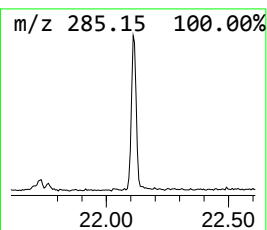
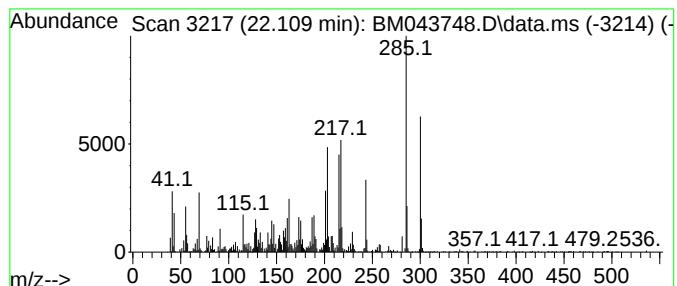
Instrument :
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Quant Title : SVOA CALIBRATION

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

Peak Number 21 9(1H)-Phenanthrenone, 2,3,4... Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.	
22.110	3.89 ng/ul	773608	Chrysene-d12	21.533	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	9(1H)-Phenanthrenone, 2,3,4,4a,1...	300	C20H28O2	000511-05-7	98
2	9(1H)-Phenanthrenone, 2,3,4,4a,1...	300	C20H28O2	000511-05-7	96
3	9(1H)-Phenanthrenone, 2,3,4,4a,1...	300	C20H28O2	000511-05-7	94
4	2(1H)-Phenanthrenone, 3,4,4a,9,1...	300	C20H28O2	006755-93-7	89
5	Phenanthrene, 1,2,3,4,4a,9,10,10...	300	C21H32O	010064-26-3	50



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM010224\
Data File : BM043748.D
Acq On : 04 Jan 2024 00:09
Operator : MA/JU
Sample : 05951-04
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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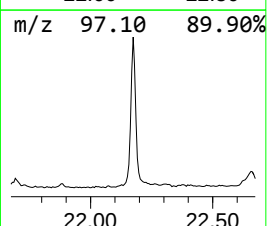
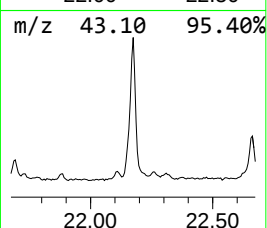
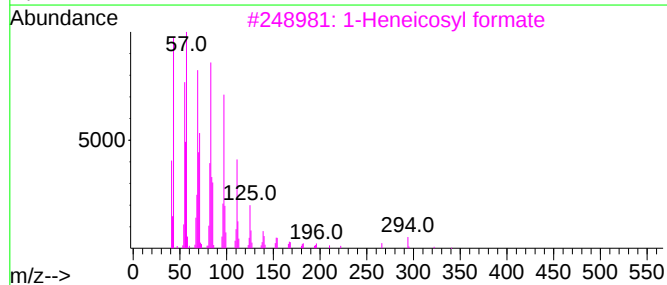
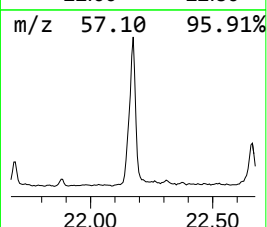
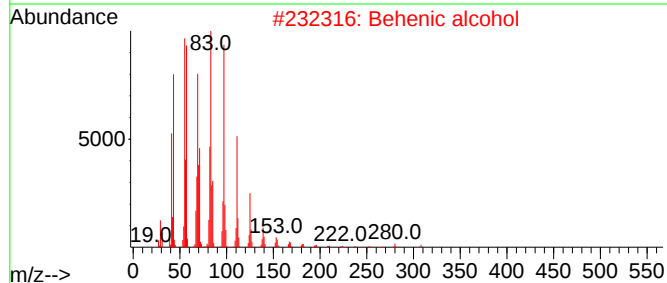
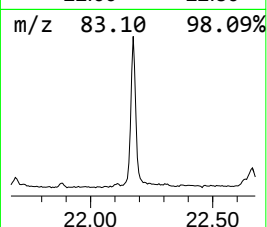
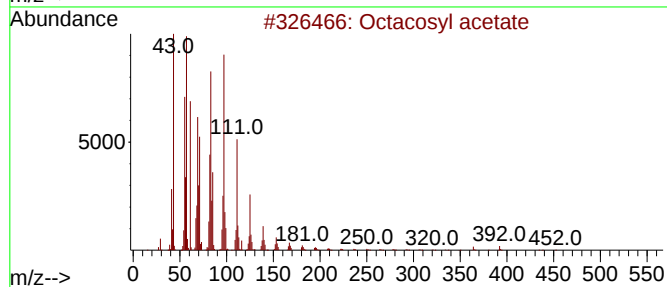
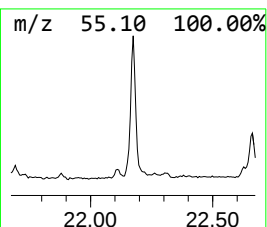
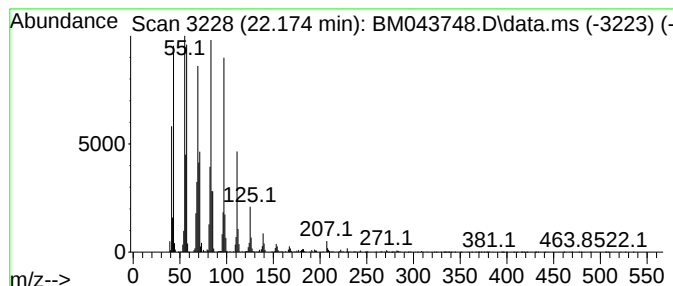
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Quant Title : SVOA CALIBRATION

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

Peak Number 22 Octacosyl acetate Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.174	20.38 ng/ul	4049890	Chrysene-d12	21.533

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octacosyl acetate	452	C30H60O2	018206-97-8	95
2			Behenic alcohol	326	C22H46O	000661-19-8	94
3			1-Heneicosyl formate	340	C22H44O2	077899-03-7	94
4			1-Tetracosene	336	C24H48	010192-32-2	94
5			n-Tetracosanol-1	354	C24H50O	000506-51-4	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM010224\
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Sample : 05951-04
Misc :
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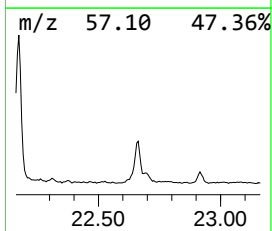
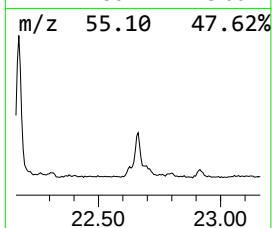
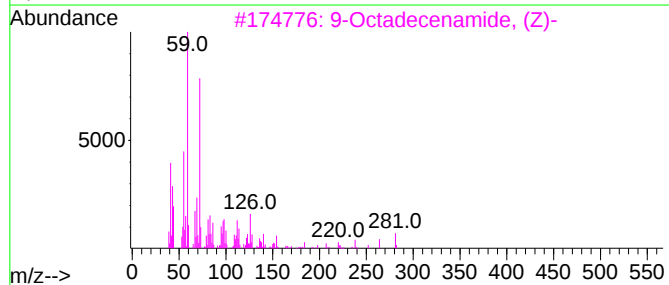
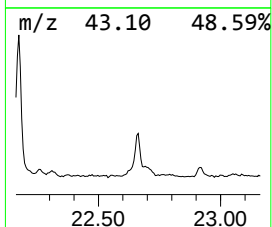
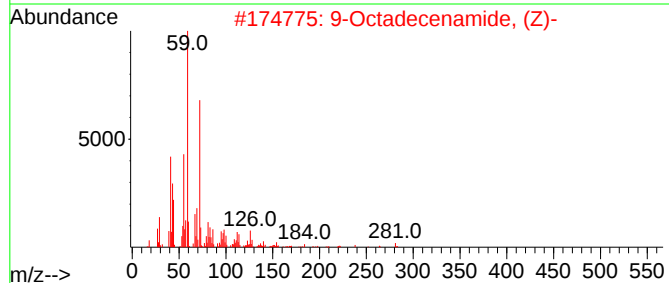
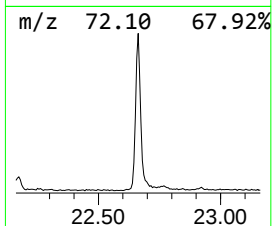
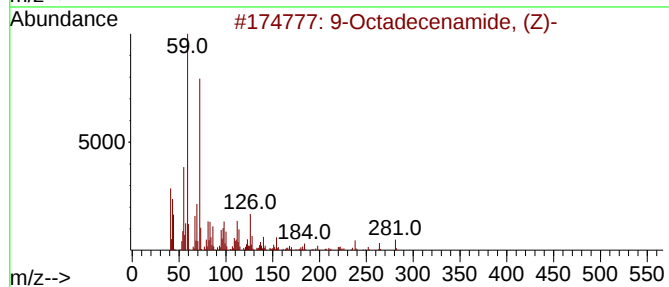
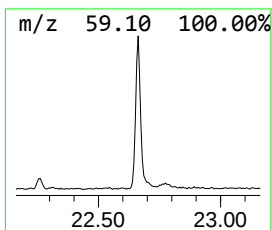
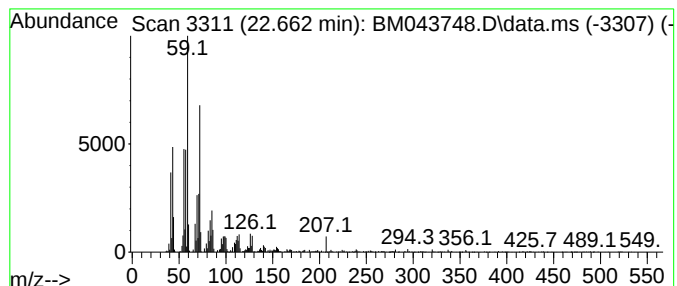
Instrument :
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ClientSampleId :
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Quant Title : SVOA CALIBRATION

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

Peak Number 23 9-Octadecenamide, (Z)- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.	
22.662	9.03 ng/ul	1794030	Chrysene-d12	21.533	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	9-Octadecenamide, (Z)-	281	C18H35NO	000301-02-0	94
2	9-Octadecenamide, (Z)-	281	C18H35NO	000301-02-0	91
3	9-Octadecenamide, (Z)-	281	C18H35NO	000301-02-0	76
4	9-Octadecenamide, (Z)-	281	C18H35NO	000301-02-0	68
5	13-Docosenamide, (Z)-	337	C22H43NO	000112-84-5	60



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM010224\
Data File : BM043748.D
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Sample : 05951-04
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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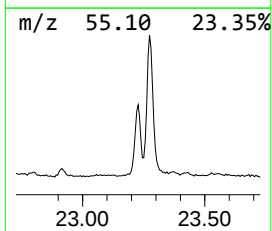
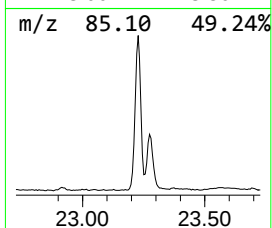
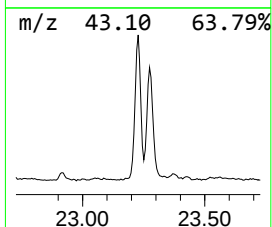
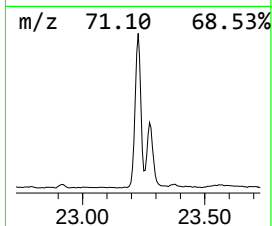
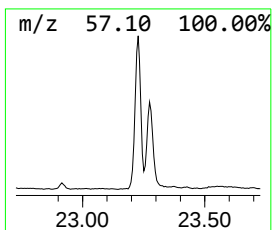
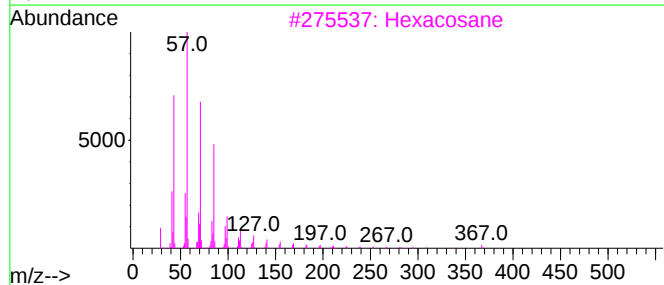
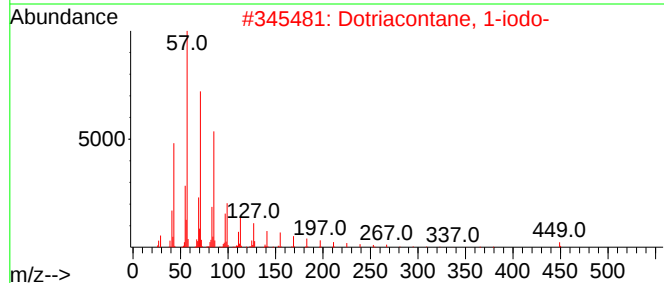
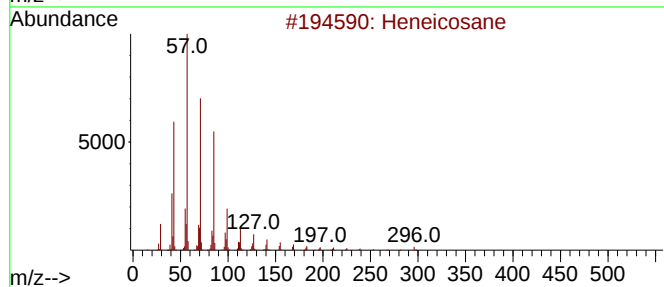
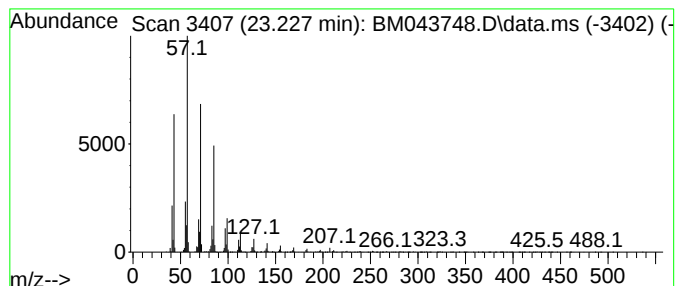
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Quant Title : SVOA CALIBRATION

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

Peak Number 24 (DEL) Alkane: Straight-Chai... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.227	16.32 ng/ul	3528600	Perylene-d12	23.956

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Heneicosane	296	C21H44	000629-94-7	91
2		Dotriacontane, 1-iodo-	576	C32H65I	1000406-32-4	91
3		Hexacosane	366	C26H54	000630-01-3	91
4		Heptadecane, 9-octyl-	352	C25H52	007225-64-1	90
5		Docosane, 11-butyl-	366	C26H54	013475-76-8	86



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM010224\
Data File : BM043748.D
Acq On : 04 Jan 2024 00:09
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Sample : 05951-04
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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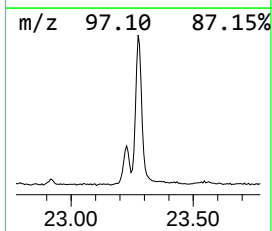
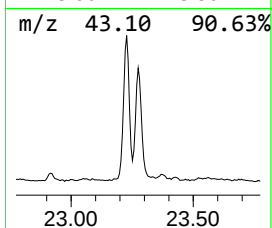
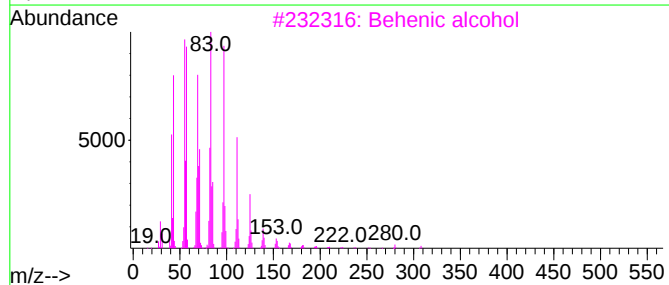
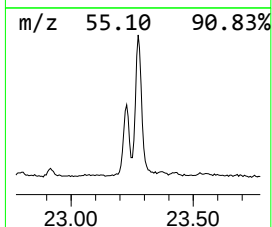
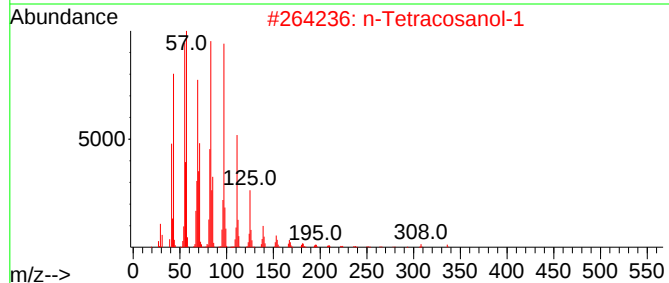
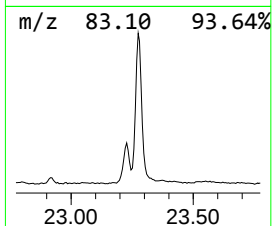
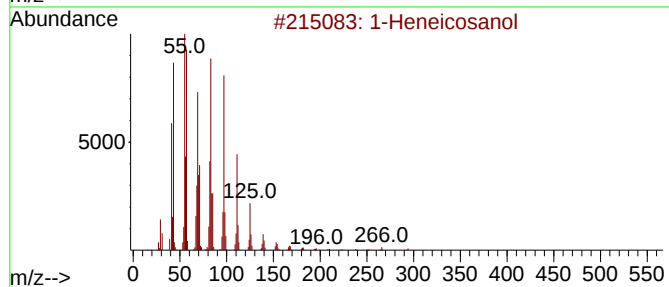
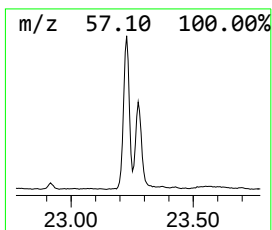
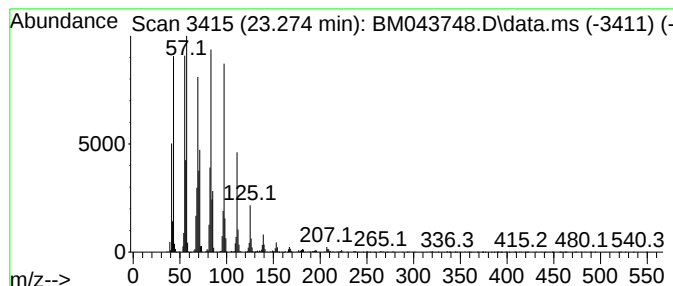
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Quant Title : SVOA CALIBRATION

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

Peak Number 25 1-Heneicosanol Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.274	21.21 ng/ul	4584940	Perylene-d12	23.956

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	94
4		Heptacos-1-ene	378	C27H54	015306-27-1	94
5		Octacosanol	410	C28H58O	000557-61-9	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM010224\
Data File : BM043748.D
Acq On : 04 Jan 2024 00:09
Operator : MA/JU
Sample : 05951-04
Misc :
ALS Vial : 16 Sample Multiplier: 1

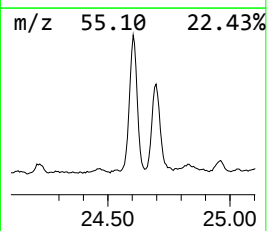
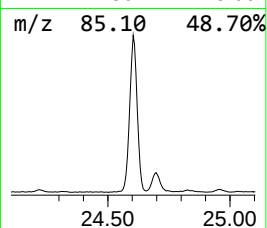
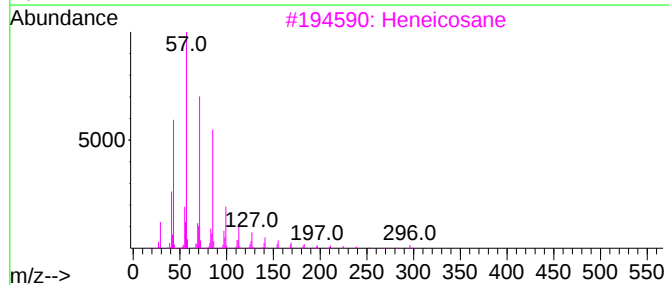
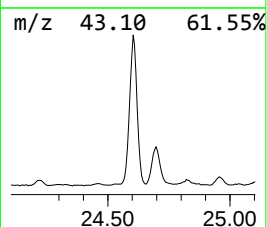
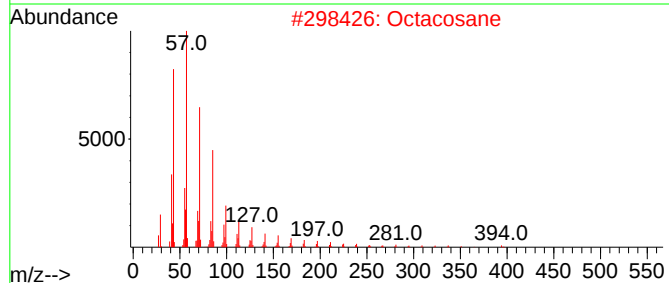
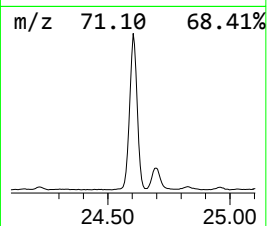
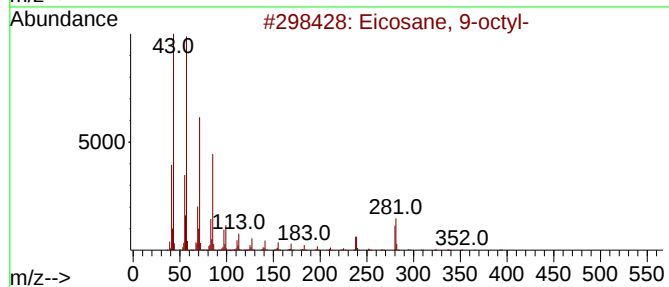
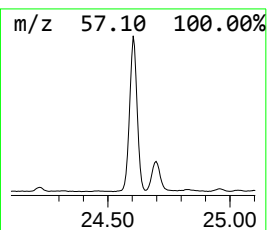
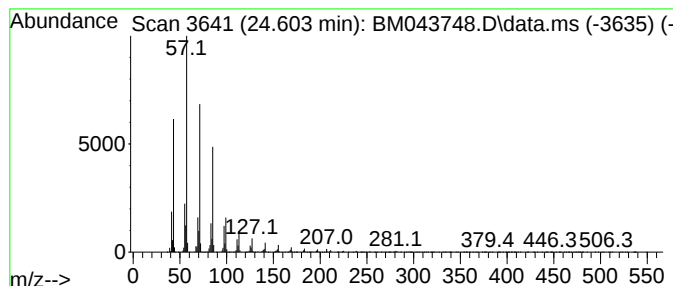
Instrument :
BNA_M
ClientSampleId :
GCF68

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 26 (DEL) Alkane: Straight-Chai... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.	
24.604	38.50 ng/ul	8322540	Perylene-d12	23.956	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Eicosane, 9-octyl-	394	C28H58	013475-77-9	94
2	Octacosane	394	C28H58	000630-02-4	91
3	Heneicosane	296	C21H44	000629-94-7	91
4	Octacosane, 2-methyl-	408	C29H60	001560-98-1	91
5	Heptadecane, 9-octyl-	352	C25H52	007225-64-1	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM010224\
Data File : BM043748.D
Acq On : 04 Jan 2024 00:09
Operator : MA/JU
Sample : 05951-04
Misc :
ALS Vial : 16 Sample Multiplier: 1

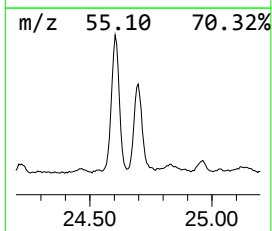
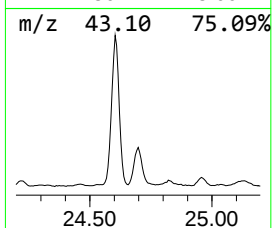
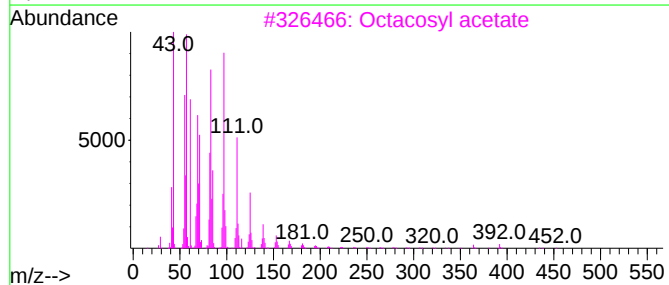
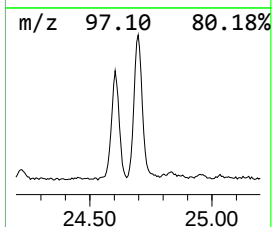
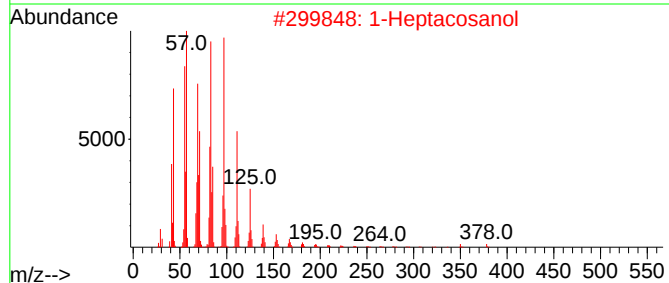
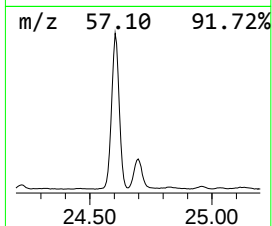
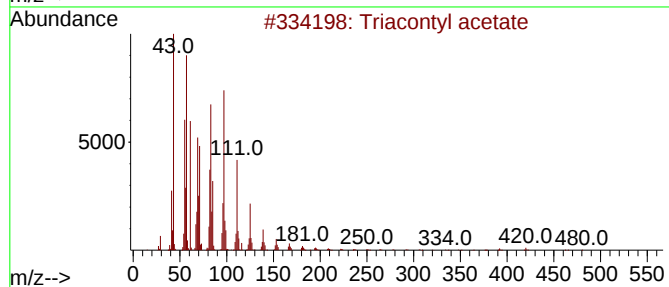
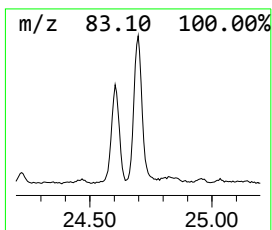
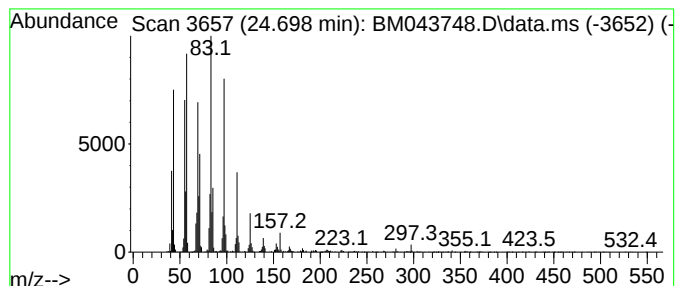
Instrument :
BNA_M
ClientSampleId :
GCF68

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 27 Triacontyl acetate Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD		R.T.	
24.698	15.45 ng/ul	3340940	Perylene-d12		23.956	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Triacontyl acetate		480	C32H64O2	041755-58-2	93
2	1-Heptacosanol		396	C27H56O	002004-39-9	93
3	Octacosyl acetate		452	C30H60O2	018206-97-8	91
4	1-Docosanol, methyl ether		340	C23H48O	1000333-91-9	90
5	1-Triacontanol		438	C30H62O	000593-50-0	86



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM010224\
Data File : BM043748.D
Acq On : 04 Jan 2024 00:09
Operator : MA/JU
Sample : 05951-04
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
GCF68

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

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Butanone	5.311	2.5	ng/ul	236870	1	7.957	1904170	20.0
Bicyclo[3.1.0]h...	6.452	2.0	ng/ul	194464	1	7.957	1904170	20.0
.alpha.-Pinene	6.616	3.1	ng/ul	295900	1	7.957	1904170	20.0
Bicyclo[3.1.0]h...	7.246	62.3	ng/ul	5927100	1	7.957	1904170	20.0
.beta.-Myrcene	7.410	4.6	ng/ul	436557	1	7.957	1904170	20.0
D-Limonene	8.163	2.8	ng/ul	268728	1	7.957	1904170	20.0
Caryophyllene	13.904	2.2	ng/ul	369804	3	14.598	3295790	20.0
(2E,4S,7E)-4-Is...	15.475	3.7	ng/ul	610560	3	14.598	3295790	20.0
n-Hexadecanoic ...	18.245	9.0	ng/ul	1725310	4	17.345	3852150	20.0
Phytol	19.257	2.2	ng/ul	430776	4	17.345	3852150	20.0
(4aS,4bR,10aS)-...	19.439	4.9	ng/ul	940704	4	17.345	3852150	20.0
(DEL) Alkane: S...	20.251	2.5	ng/ul	494562	5	21.533	3975300	20.0
(1R,4aR,4bS,10a...	20.286	4.1	ng/ul	806808	5	21.533	3975300	20.0
2-Phenanthrenol...	20.445	3.7	ng/ul	734005	5	21.533	3975300	20.0
Cycloisolongifo...	20.616	29.4	ng/ul	5850030	5	21.533	3975300	20.0
unknown-01	20.657	10.4	ng/ul	2065310	5	21.533	3975300	20.0
(1R,4aR,4bS,10a...	20.980	7.9	ng/ul	1561580	5	21.533	3975300	20.0
((1R,4aR,4bR,10...	21.115	17.0	ng/ul	3374920	5	21.533	3975300	20.0
(DEL) Alkane: S...	21.239	12.9	ng/ul	2553330	5	21.533	3975300	20.0
unknown-02	21.339	4.8	ng/ul	948114	5	21.533	3975300	20.0
9(1H)-Phenanthr...	22.110	3.9	ng/ul	773608	5	21.533	3975300	20.0
Octacosyl acetate	22.174	20.4	ng/ul	4049890	5	21.533	3975300	20.0
9-Octadecenamid...	22.662	9.0	ng/ul	1794030	5	21.533	3975300	20.0
(DEL) Alkane: S...	23.227	16.3	ng/ul	3528600	6	23.956	4323710	20.0
1-Heneicosanol	23.274	21.2	ng/ul	4584940	6	23.956	4323710	20.0
(DEL) Alkane: S...	24.604	38.5	ng/ul	8322540	6	23.956	4323710	20.0
Triacetyl acetate	24.698	15.4	ng/ul	3340940	6	23.956	4323710	20.0