

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN060624\
 Data File : BN031820.D
 Acq On : 06 Jun 2024 15:55
 Operator : MA/JU
 Sample : P2634-06
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 BHBX0

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_N\Methods\SFAM-EPA-BN052924.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BN031820.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.181	30	33	45	rVB	84820	123252	1.35%	0.088%
2	3.446	74	78	93	rVB	218147	314747	3.46%	0.224%
3	3.587	99	102	116	rBV	1055675	1541383	16.94%	1.098%
4	5.769	468	473	483	rVB	63277	101811	1.12%	0.073%
5	6.358	567	573	582	rVB	659946	1029256	11.31%	0.733%
6	6.663	620	625	631	rVB	231581	384511	4.23%	0.274%
7	6.846	643	656	671	rVB	1508302	2469267	27.14%	1.759%
8	7.011	671	684	694	rBV	1212062	1969453	21.65%	1.403%
9	7.105	694	700	710	rVB	1034593	1757056	19.31%	1.251%
10	7.205	710	717	725	rBV	1511798	2418461	26.58%	1.723%
11	7.281	725	730	736	rBV	43210	74238	0.82%	0.053%
12	7.563	772	778	787	rBV	145621	235276	2.59%	0.168%
13	7.669	789	796	803	rVV	1461115	2319726	25.50%	1.652%
14	7.740	803	808	814	rVV2	55821	102051	1.12%	0.073%
15	7.887	827	833	839	rVV	156649	247282	2.72%	0.176%
16	8.375	909	916	925	rBV	1900671	3050576	33.53%	2.173%
17	8.828	985	993	1000	rBV	1398097	2285083	25.12%	1.628%
18	9.393	1082	1089	1096	rBV	189448	293081	3.22%	0.209%
19	9.546	1107	1115	1127	rBV	1102835	1870130	20.55%	1.332%
20	10.075	1197	1205	1227	rVB	1751356	3013300	33.12%	2.146%
21	10.452	1260	1269	1277	rBV	1960396	3370938	37.05%	2.401%
22	10.593	1283	1293	1312	rBV	1569992	2774758	30.50%	1.976%
23	10.999	1356	1362	1370	rBV2	73059	132615	1.46%	0.094%
24	11.199	1389	1396	1416	rVB	192470	408580	4.49%	0.291%
25	11.822	1495	1502	1508	rBV	87978	141490	1.56%	0.101%
26	13.187	1729	1734	1738	rVV2	52531	90538	1.00%	0.064%
27	13.251	1740	1745	1753	rVV4	33822	64606	0.71%	0.046%
28	13.634	1804	1810	1815	rVV2	77171	123192	1.35%	0.088%
29	13.728	1820	1826	1832	rVV	2708300	4007557	44.05%	2.854%
30	13.804	1836	1839	1842	rVV3	46708	69960	0.77%	0.050%
31	13.845	1842	1846	1854	rVB2	165065	254686	2.80%	0.181%
32	14.010	1867	1874	1890	rVB	3293473	5357294	58.88%	3.816%
33	14.275	1913	1919	1921	rVV	157527	235675	2.59%	0.168%
34	14.316	1921	1926	1932	rVB	2863204	4297303	47.23%	3.061%
35	14.387	1932	1938	1951	rVB5	99919	253233	2.78%	0.180%
36	14.522	1955	1961	1969	rBV	1353391	2176473	23.92%	1.550%

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Integration Parameters: LSCINT.P

Integrator: RTE

Smoothing : OFF

Filtering: 5

Sampling : 1

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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_N\Methods\SFAM-EPA-BN052924.MA.M
 Title : SVOA CALIBRATION

37	14.592	1969	1973	1981	rVB2	144006	229269	2.52%	0.163%
38	14.869	2014	2020	2025	rVB	404932	552678	6.07%	0.394%
39	15.310	2088	2095	2101	rBV	4305503	6160957	67.72%	4.388%
40	15.369	2101	2105	2111	rVB4	49372	83451	0.92%	0.059%
41	15.440	2111	2117	2129	rBV	1085777	1567357	17.23%	1.116%
42	15.657	2149	2154	2160	rBV3	40594	73617	0.81%	0.052%
43	15.798	2171	2178	2183	rBV4	62237	103924	1.14%	0.074%
44	15.939	2197	2202	2213	rVB2	81111	126996	1.40%	0.090%
45	16.039	2214	2219	2225	rBV4	56176	99940	1.10%	0.071%
46	16.722	2330	2335	2342	rVB2	50517	92697	1.02%	0.066%
47	16.881	2359	2362	2370	rVB	55144	86328	0.95%	0.061%
48	17.063	2386	2393	2397	rBV	3455841	5041659	55.41%	3.591%
49	17.104	2397	2400	2404	rVV	917921	1271994	13.98%	0.906%
50	17.163	2404	2410	2421	rVV	4409112	6463072	71.04%	4.603%
51	17.251	2421	2425	2430	rVV5	42400	82136	0.90%	0.059%
52	17.469	2454	2462	2469	rBV	82311	165721	1.82%	0.118%
53	17.639	2488	2491	2498	rVB3	59273	99149	1.09%	0.071%
54	17.939	2537	2542	2543	rBV	207270	254817	2.80%	0.181%
55	17.969	2543	2547	2556	rVB2	554280	1224936	13.46%	0.872%
56	18.069	2561	2564	2569	rVV	66100	110144	1.21%	0.078%
57	18.128	2569	2574	2578	rVV2	289265	523200	5.75%	0.373%
58	18.169	2578	2581	2586	rVB	160289	228357	2.51%	0.163%
59	18.445	2623	2628	2631	rBV	167709	231410	2.54%	0.165%
60	18.486	2631	2635	2645	rVB2	166055	263048	2.89%	0.187%
61	18.739	2675	2678	2681	rVV	73450	88803	0.98%	0.063%
62	18.781	2682	2685	2689	rVB	61507	75044	0.82%	0.053%
63	18.863	2694	2699	2703	rBV	197189	328658	3.61%	0.234%
64	18.922	2704	2709	2712	rBV4	64615	103655	1.14%	0.074%
65	18.951	2712	2714	2718	rVB	58953	68581	0.75%	0.049%
66	19.004	2719	2723	2731	rBV4	127750	305298	3.36%	0.217%
67	19.128	2735	2744	2751	rBV	2259302	3240709	35.62%	2.308%
68	19.251	2758	2765	2768	rBV4	186552	336669	3.70%	0.240%
69	19.463	2795	2801	2804	rBV	5628654	8208694	90.22%	5.847%
70	19.492	2804	2806	2814	rVV	2088626	2480741	27.27%	1.767%
71	19.563	2815	2818	2824	rVB2	119441	197151	2.17%	0.140%
72	19.675	2833	2837	2842	rVB	72204	115529	1.27%	0.082%
73	19.733	2844	2847	2852	rVB	88651	129962	1.43%	0.093%
74	19.816	2856	2861	2867	rBV2	176451	439245	4.83%	0.313%
75	19.951	2880	2884	2887	rBV6	124062	231633	2.55%	0.165%
76	19.986	2887	2890	2894	rVB2	296514	383063	4.21%	0.273%

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Integration Parameters: LSCINT.P

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

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 Peak separation: 5

Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN052924.MA.M
 Title : SVOA CALIBRATION

77	20.086	2904	2907	2911	rBV	176888	237046	2.61%	0.169%
78	20.145	2913	2917	2921	rVB2	230124	336903	3.70%	0.240%
79	20.286	2938	2941	2945	rVB2	164173	183221	2.01%	0.130%
80	20.739	3015	3018	3023	rVB	194926	246804	2.71%	0.176%
81	20.951	3051	3054	3056	rBV	169639	189361	2.08%	0.135%
82	20.986	3056	3060	3067	rVB2	911319	1260160	13.85%	0.898%
83	21.210	3096	3098	3102	rVB	162391	166473	1.83%	0.119%
84	21.298	3106	3113	3118	rVV3	3504224	6427307	70.64%	4.578%
85	21.345	3118	3121	3130	rVB	944116	1349215	14.83%	0.961%
86	21.992	3226	3231	3237	rVV2	2087799	3721666	40.91%	2.651%
87	22.051	3238	3241	3249	rVB7	443407	798009	8.77%	0.568%
88	22.157	3251	3259	3265	rVB4	1934204	3883756	42.69%	2.766%
89	22.245	3271	3274	3277	rBV3	265954	392226	4.31%	0.279%
90	22.292	3277	3282	3286	rVV	1283303	2127294	23.38%	1.515%
91	22.339	3287	3290	3297	rVB2	506029	781082	8.59%	0.556%
92	22.451	3303	3309	3318	rVB2	1238278	2863345	31.47%	2.039%
93	22.610	3331	3336	3342	rBV4	977968	2119941	23.30%	1.510%
94	22.669	3342	3346	3351	rVB3	680536	1124066	12.35%	0.801%
95	22.839	3370	3375	3387	rVB	565867	1080937	11.88%	0.770%
96	23.057	3409	3412	3417	rVB2	408789	576601	6.34%	0.411%
97	23.145	3417	3427	3438	rVB4	3633585	9098182	100.00%	6.480%
98	23.316	3448	3456	3463	rBV3	831672	2347007	25.80%	1.672%
99	24.039	3572	3579	3585	rBV2	1892476	4220287	46.39%	3.006%
100	24.239	3605	3613	3622	rVB	1506925	3641038	40.02%	2.593%

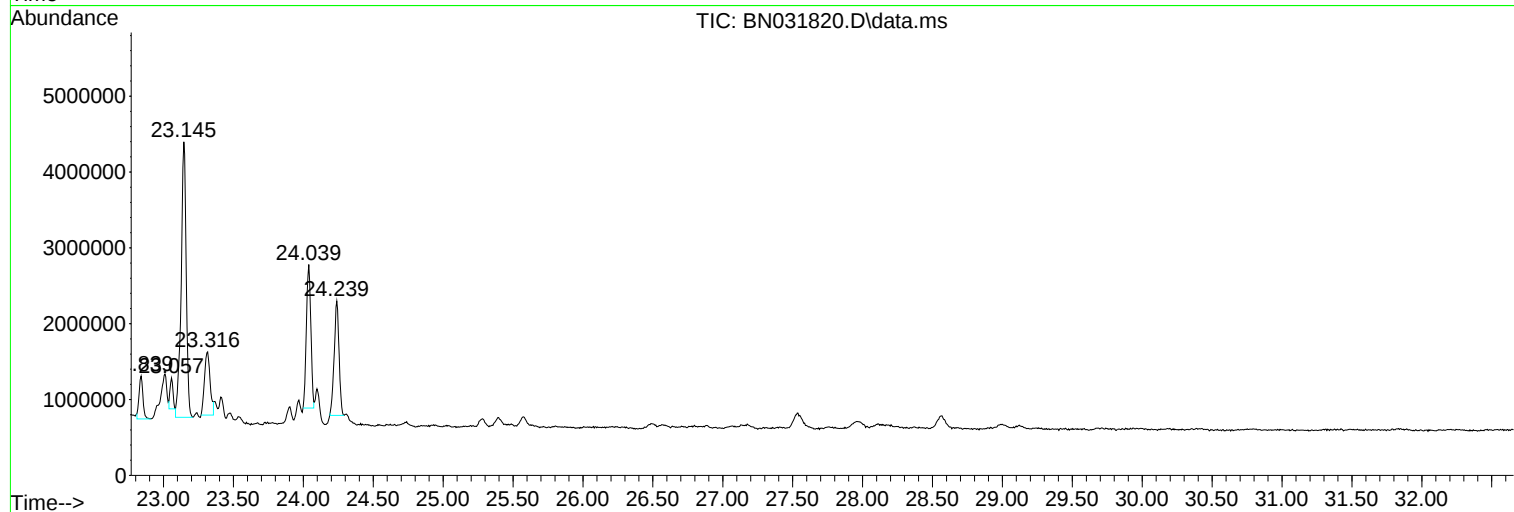
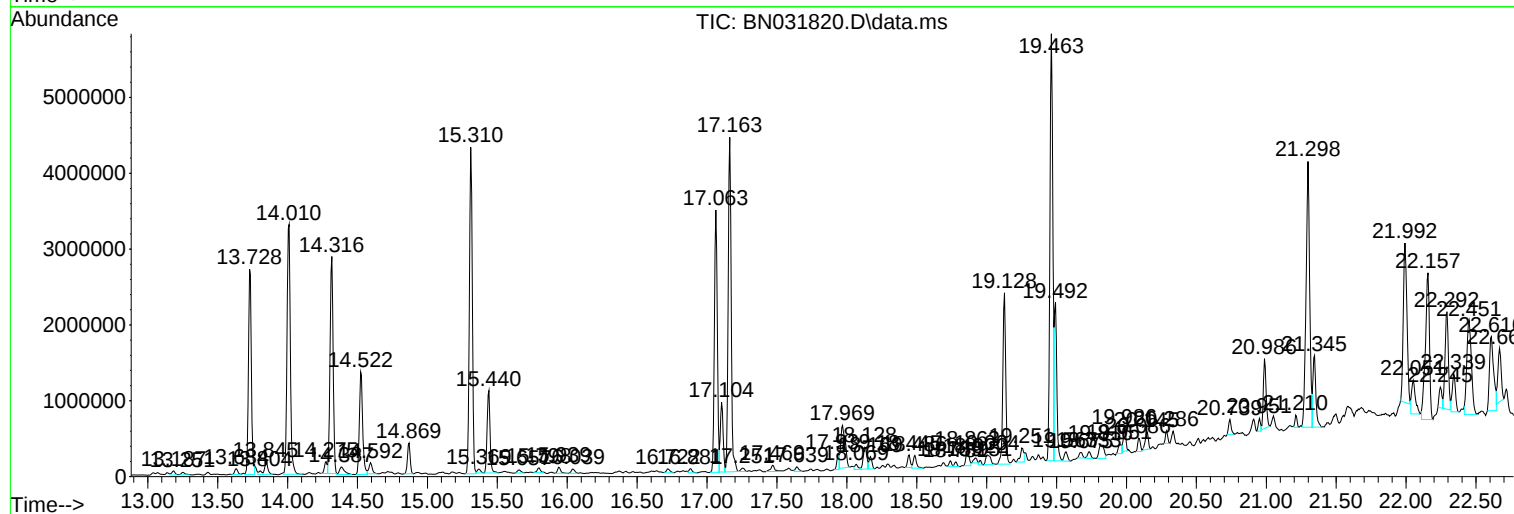
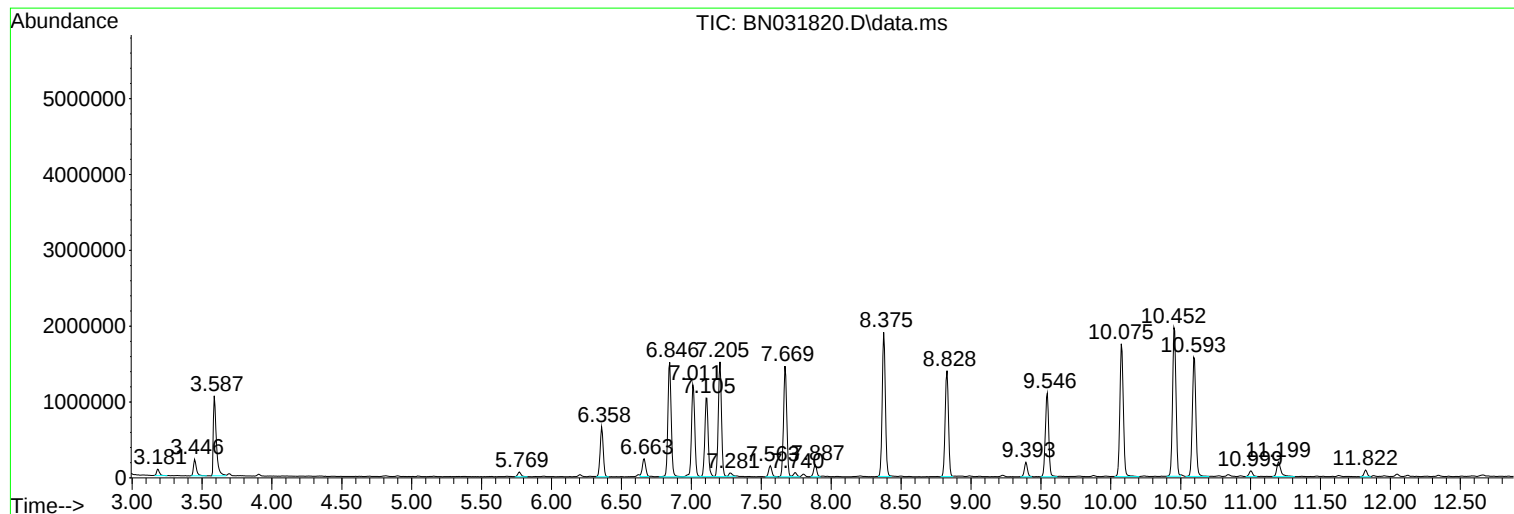
Sum of corrected areas: 140401057

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN052924.MA.M
Quant Title : SVOA CALIBRATION

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



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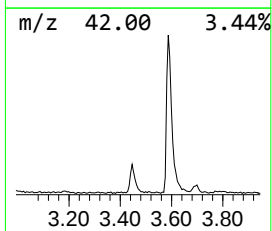
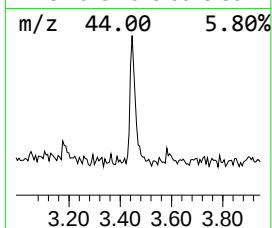
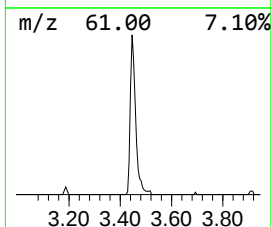
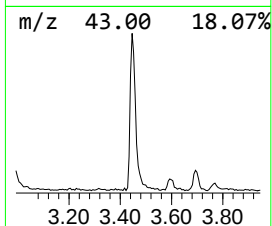
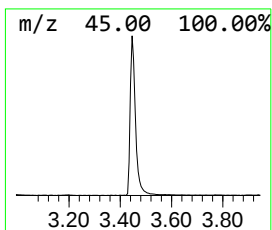
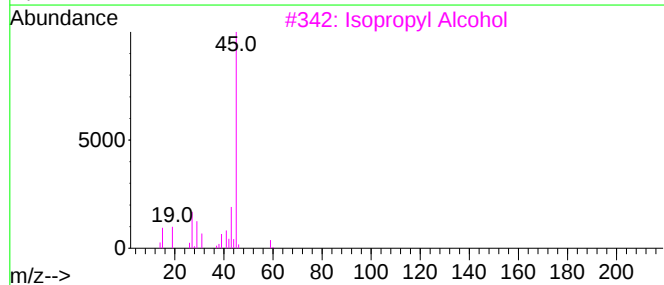
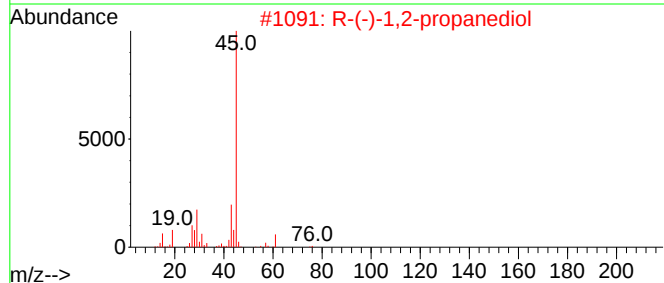
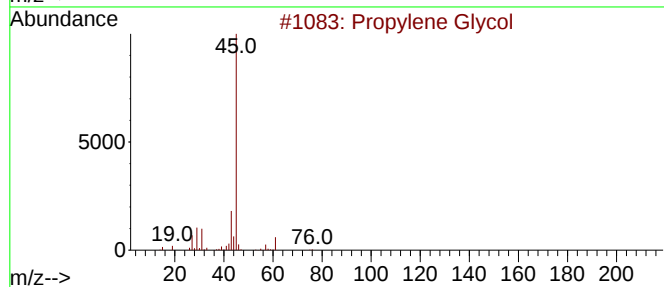
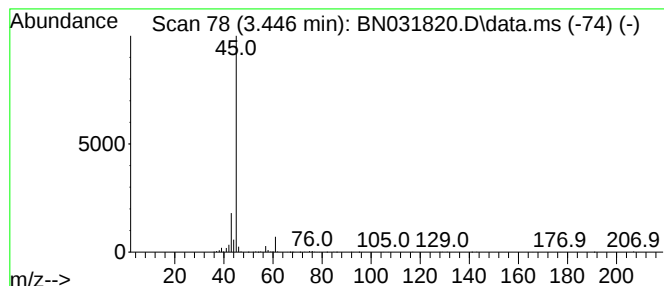
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN052924.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Propylene Glycol Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.446	2.71 ng/ul	314747	1,4-Dichlorobenzene-d4	7.669

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Propylene Glycol	76	C3H8O2	000057-55-6	86
2			R-(-)-1,2-propanediol	76	C3H8O2	004254-14-2	78
3			Isopropyl Alcohol	60	C3H8O	000067-63-0	56
4			Propylene Glycol	76	C3H8O2	000057-55-6	56
5			(S)-(+)-1,2-Propanediol	76	C3H8O2	004254-15-3	40



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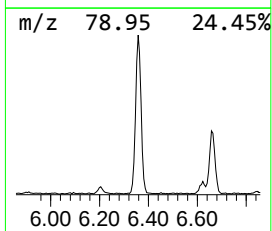
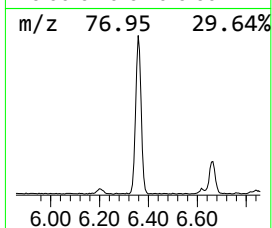
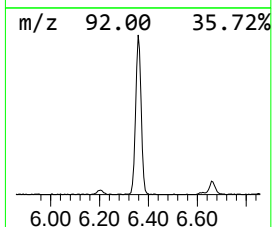
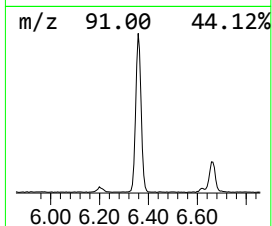
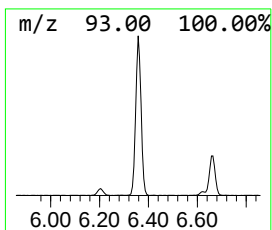
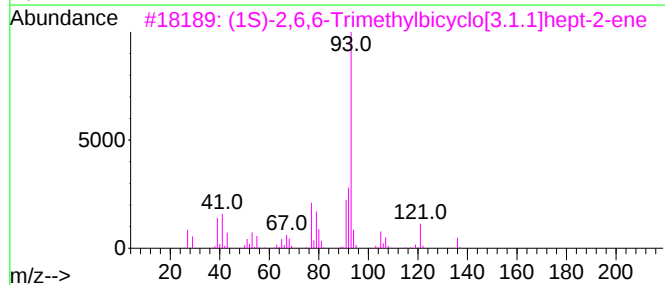
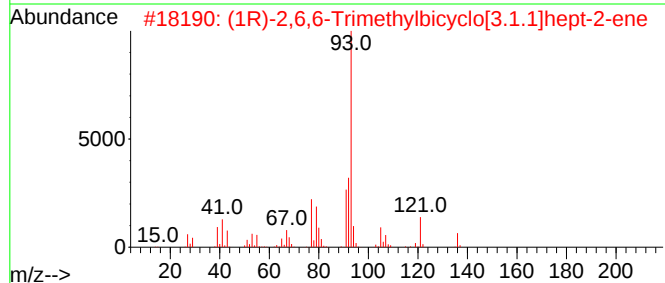
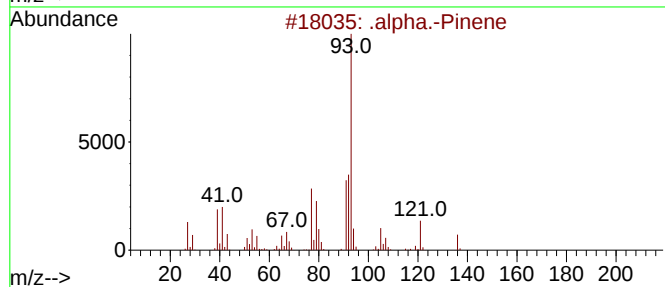
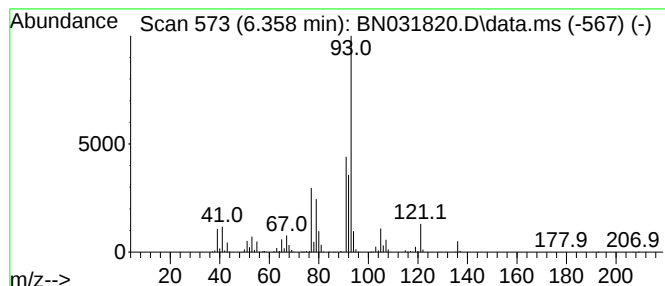
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN052924.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 .alpha.-Pinene Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.358	8.87 ng/ul	1029260	1,4-Dichlorobenzene-d4	7.669

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...			136	C10H16	007785-70-8	95
3	(1S)-2,6,6-Trimethylbicyclo[3.1...			136	C10H16	007785-26-4	94
4	Bicyclo[3.1.1]hept-2-ene, 3,6,6-...			136	C10H16	004889-83-2	94
5	.		.alpha.-Pinene	136	C10H16	000080-56-8	94



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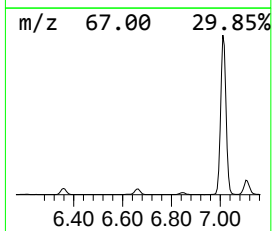
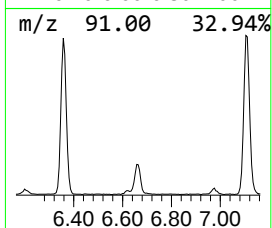
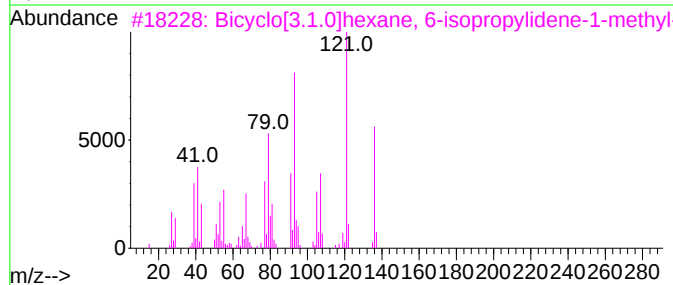
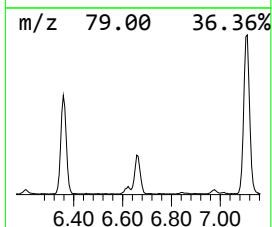
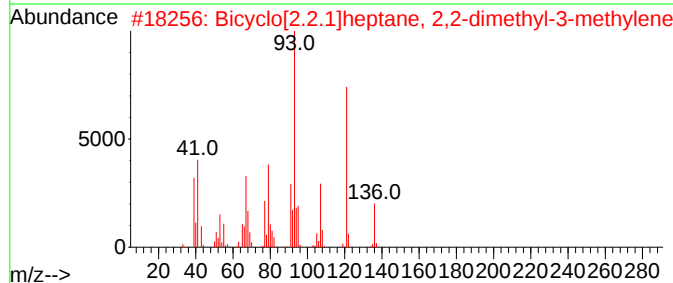
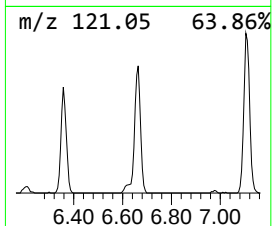
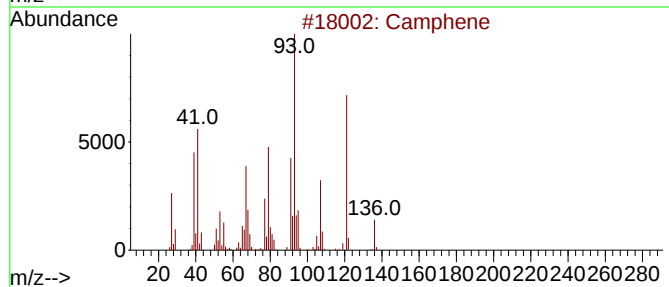
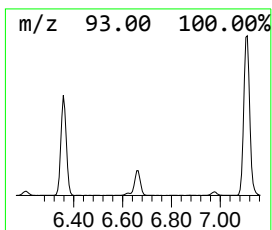
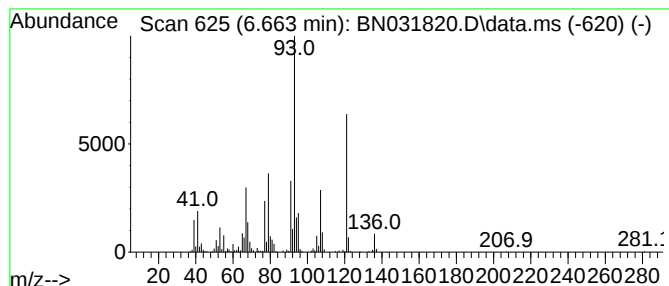
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Quant Title : SVOA CALIBRATION

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

Peak Number 3 Camphene Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.663	3.32 ng/ul	384511	1,4-Dichlorobenzene-d4	7.669

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Camphene	136	C10H16	000079-92-5	93
2			Bicyclo[2.2.1]heptane, 2,2-dimet...	136	C10H16	005794-04-7	83
3			Bicyclo[3.1.0]hexane, 6-isopropy...	136	C10H16	024524-57-0	64
4			.beta.-Pinene	136	C10H16	000127-91-3	64
5			(+)-4-Carene	136	C10H16	029050-33-7	62



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN060624\
Data File : BN031820.D
Acq On : 06 Jun 2024 15:55
Operator : MA/JU
Sample : P2634-06
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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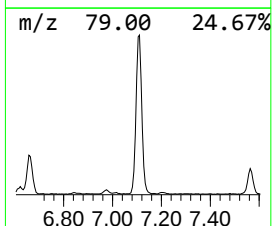
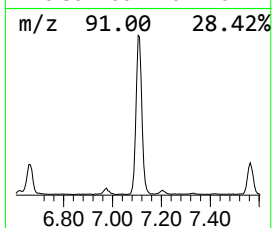
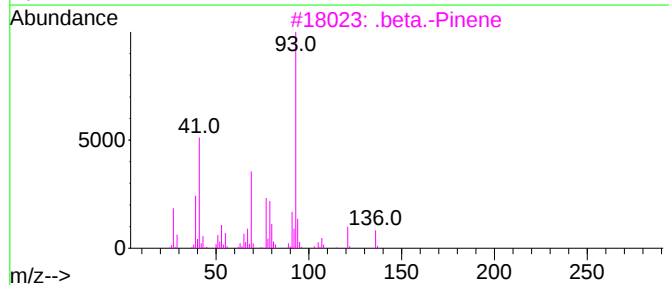
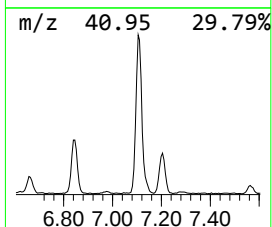
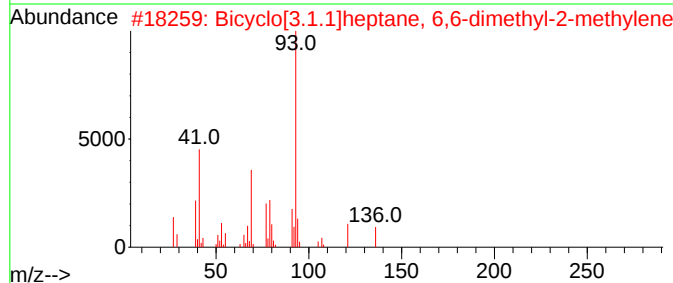
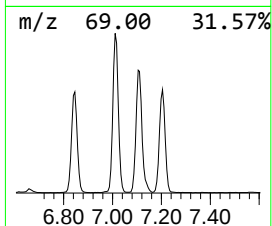
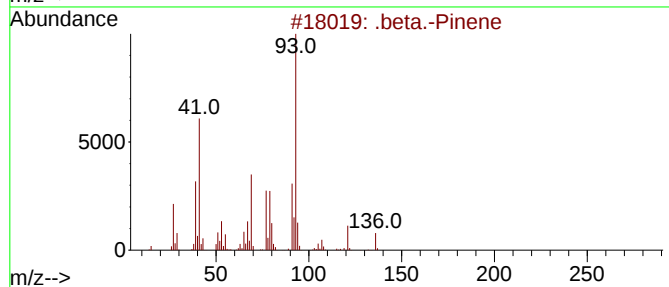
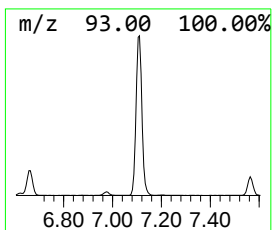
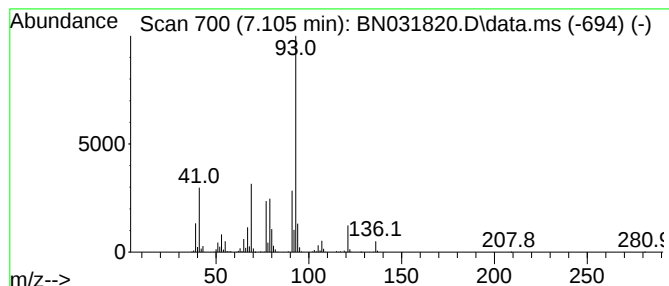
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 .beta.-Pinene Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.105	15.15 ng/ul	1757060	1,4-Dichlorobenzene-d4	7.669

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	.	beta.-Pinene	136	C10H16	000127-91-3	94	
2	Bicyclo[3.1.1]heptane, 6,6-dimet...	136	C10H16	018172-67-3	91		
3	.	beta.-Pinene	136	C10H16	000127-91-3	91	
4	Cyclohexene, 4-methylene-1-(1-me...	136	C10H16	000099-84-3	91		
5	Bicyclo[3.1.1]heptane, 6,6-dimet...	136	C10H16	018172-67-3	91		



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN060624\
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Sample : P2634-06
Misc :
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Instrument :
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ClientSampleId :
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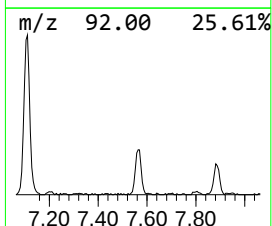
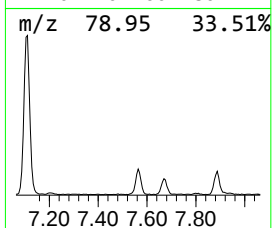
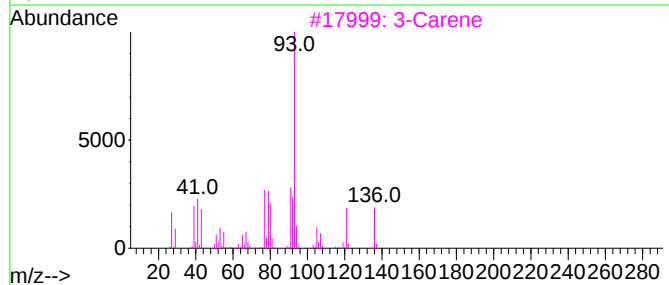
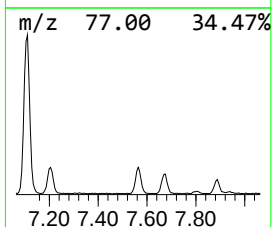
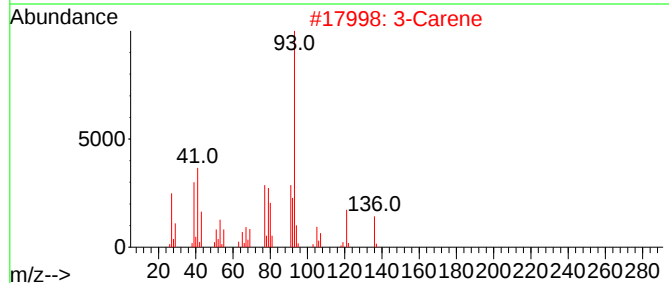
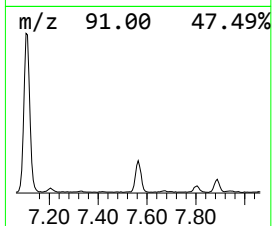
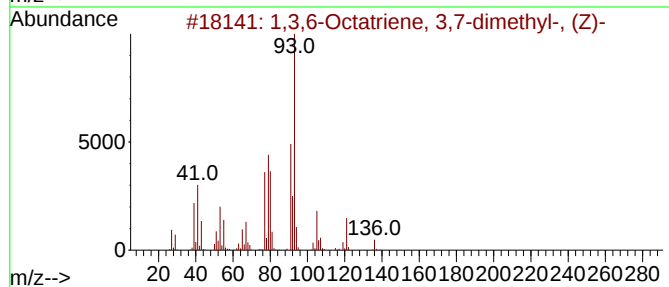
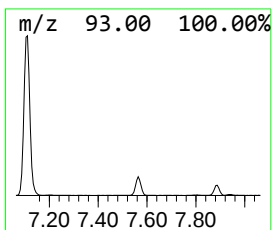
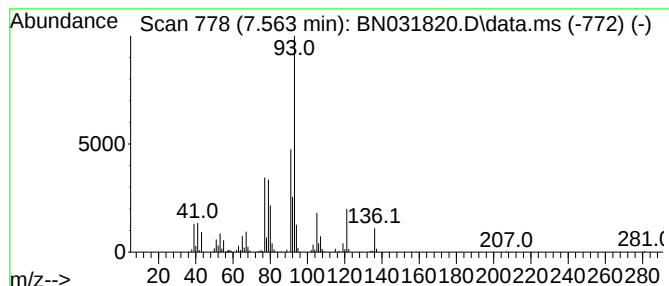
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 1,3,6-Octatriene, 3,7-dimet... Concentration Rank 22

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.563	2.03 ng/ul	235276	1,4-Dichlorobenzene-d4	7.669

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,3,6-Octatriene, 3,7-dimethyl-,...	136	C10H16	003338-55-4	96		
2	3-Carene	136	C10H16	013466-78-9	95		
3	3-Carene	136	C10H16	013466-78-9	95		
4	.beta.-Ocimene	136	C10H16	013877-91-3	94		
5	(+)-3-Carene	136	C10H16	000498-15-7	91		



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Data File : BN031820.D
Acq On : 06 Jun 2024 15:55
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Sample : P2634-06
Misc :
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Instrument :
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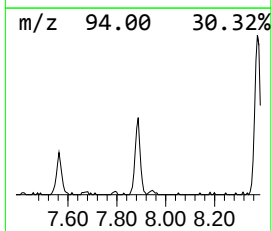
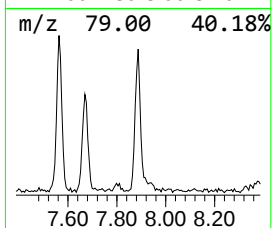
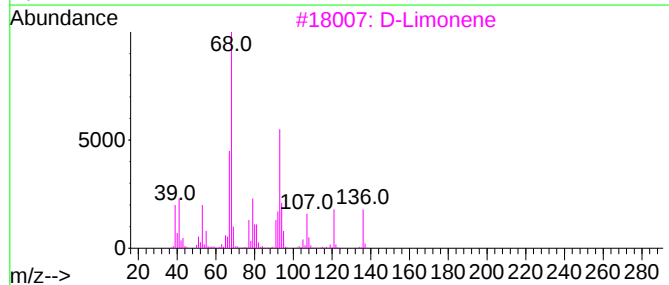
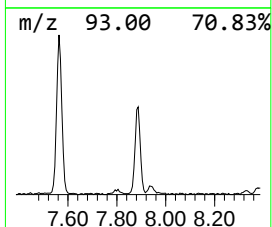
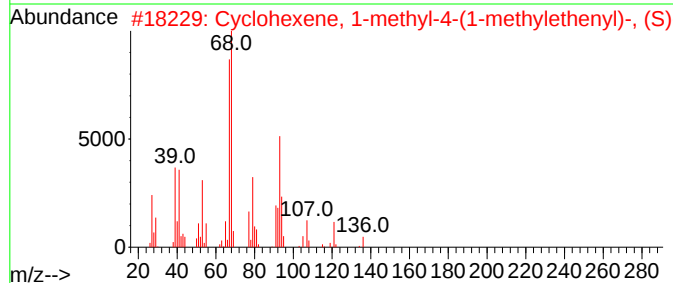
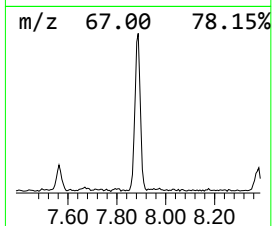
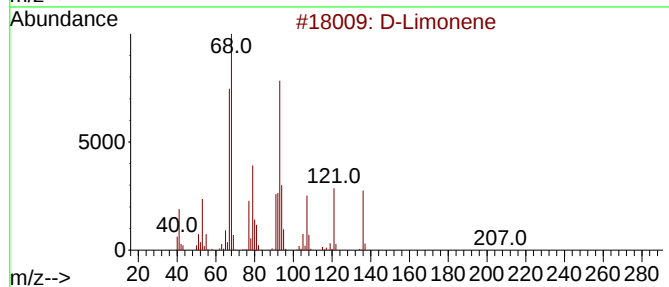
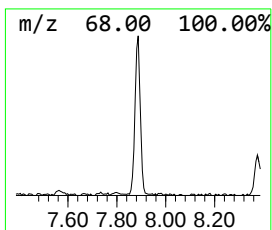
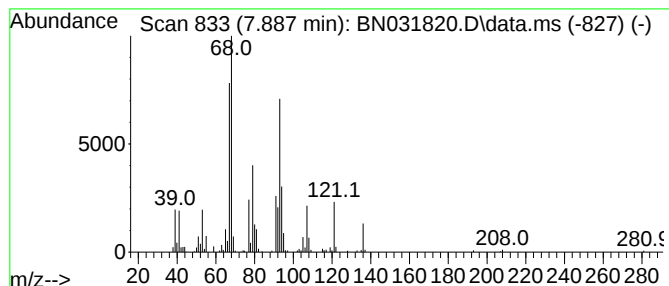
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 D-Limonene Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.887	2.13 ng/ul	247282	1,4-Dichlorobenzene-d4	7.669

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			D-Limonene	136	C10H16	005989-27-5	97
2			Cyclohexene, 1-methyl-4-(1-methy...	136	C10H16	005989-54-8	95
3			D-Limonene	136	C10H16	005989-27-5	93
4			Limonene	136	C10H16	000138-86-3	92
5			1,5-Cyclooctadiene, 1,5-dimethyl-	136	C10H16	003760-14-3	86



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN060624\
Data File : BN031820.D
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Sample : P2634-06
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
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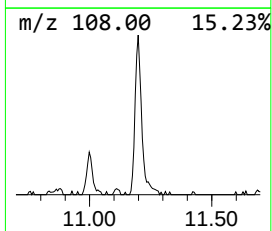
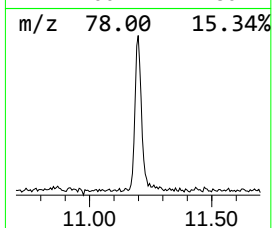
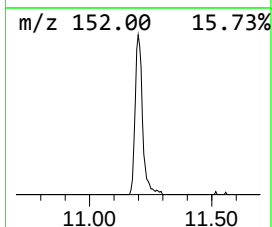
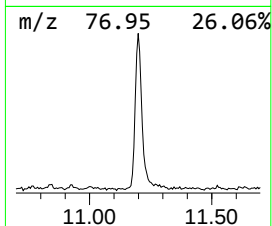
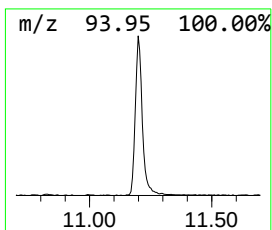
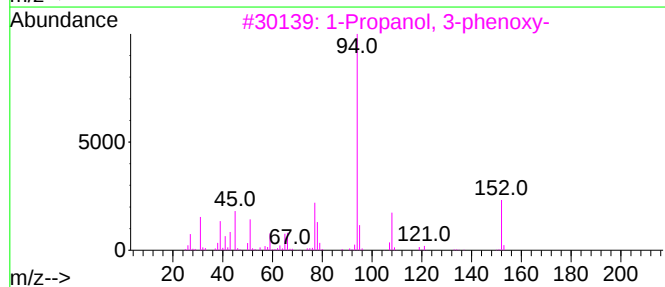
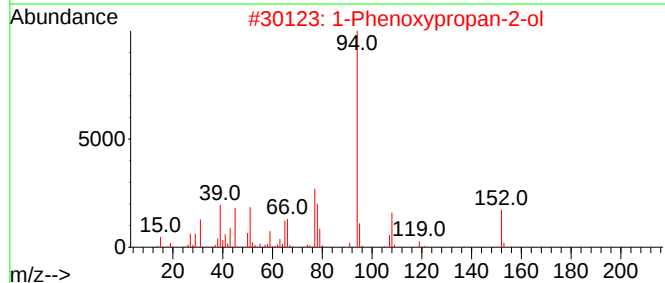
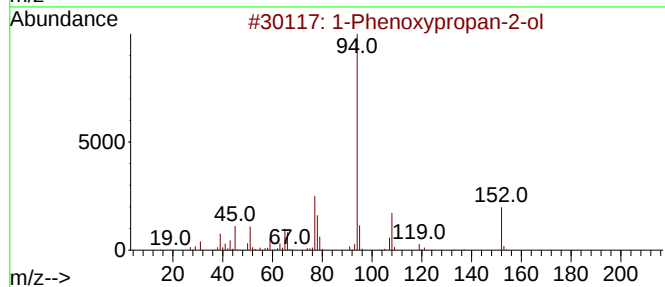
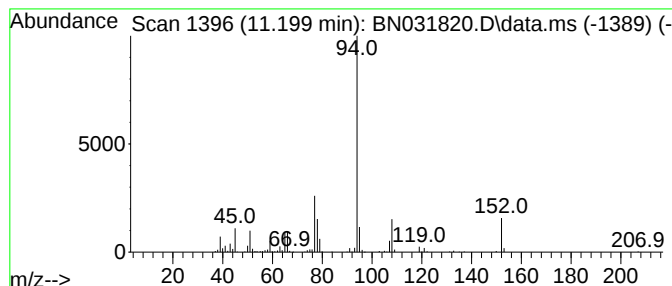
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 1-Phenoxypropan-2-ol Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.199	2.42 ng/ul	408580	Naphthalene-d8	10.451

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	96
2			1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	95
3			1-Propanol, 3-phenoxy-	152	C9H12O2	006180-61-6	91
4			1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	87
5			1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	87



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Data File : BN031820.D
Acq On : 06 Jun 2024 15:55
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Sample : P2634-06
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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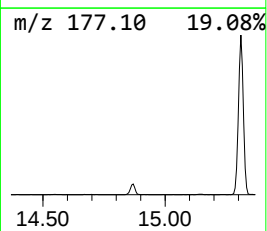
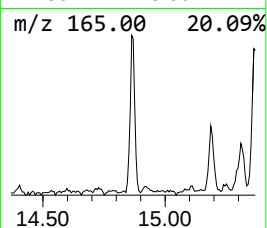
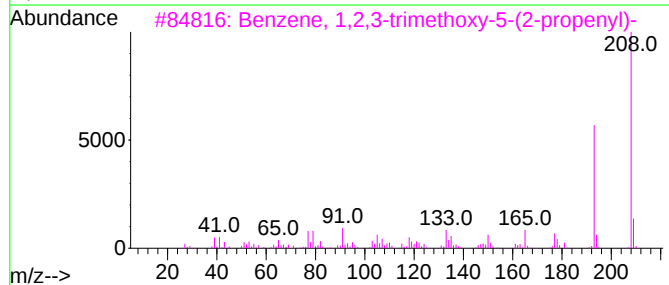
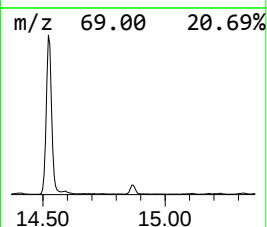
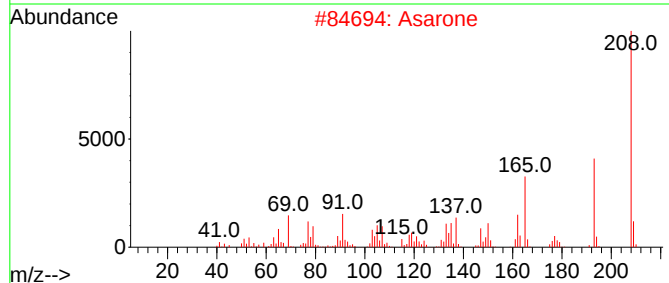
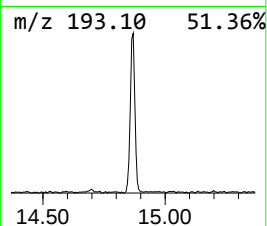
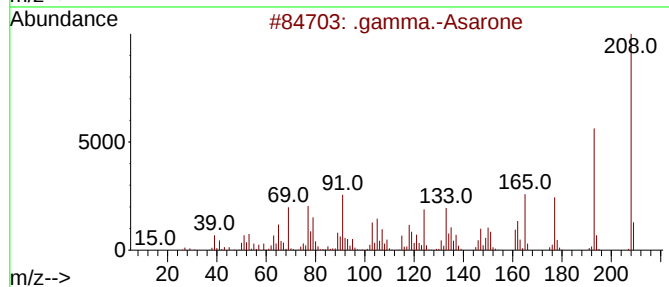
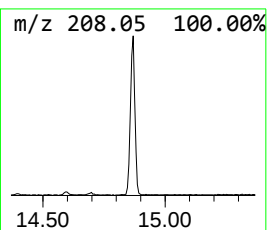
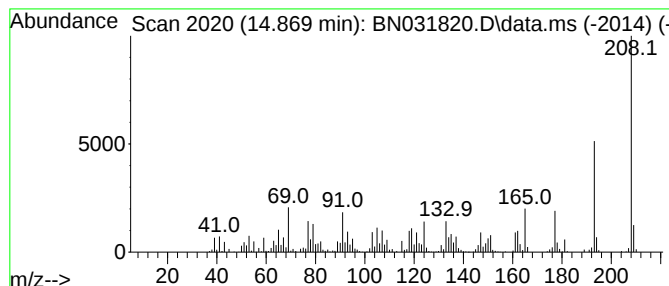
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TIC Integration Parameters: LSCINT.P

Peak Number 8 .gamma.-Asarone Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.869	2.57 ng/ul	552678	Acenaphthene-d10	14.316

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			.gamma.-Asarone	208	C12H16O3	005353-15-1	98
2			Asarone	208	C12H16O3	002883-98-9	93
3			Benzene, 1,2,3-trimethoxy-5-(2-p...	208	C12H16O3	000487-11-6	93
4			Benzene, 1,2,3-trimethoxy-5-(2-p...	208	C12H16O3	000487-11-6	92
5			.beta.-Asarone	208	C12H16O3	005273-86-9	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN060624\
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ALS Vial : 9 Sample Multiplier: 1

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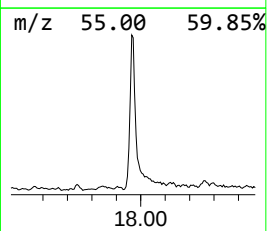
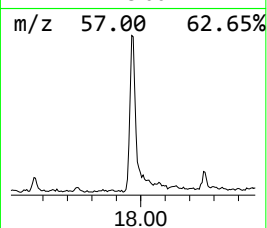
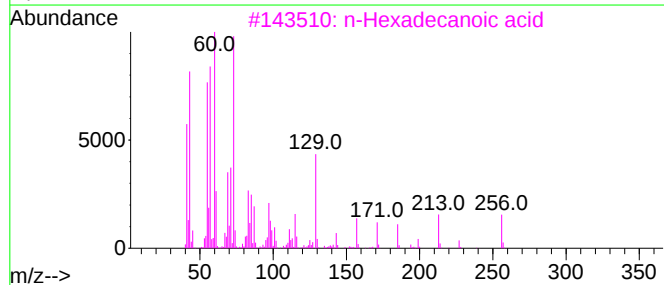
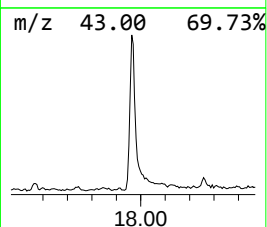
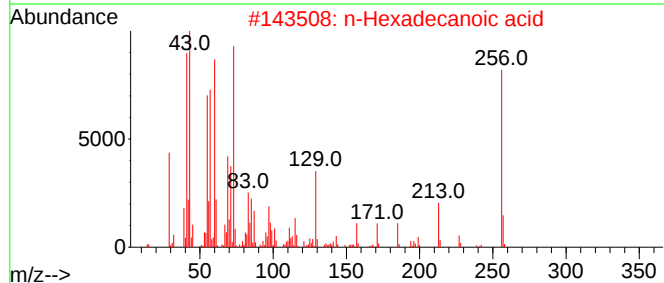
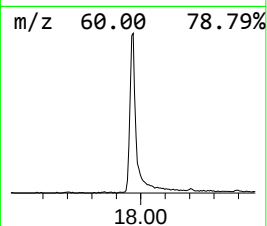
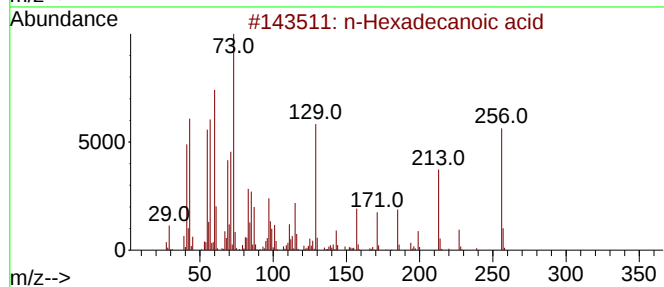
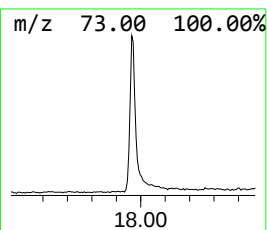
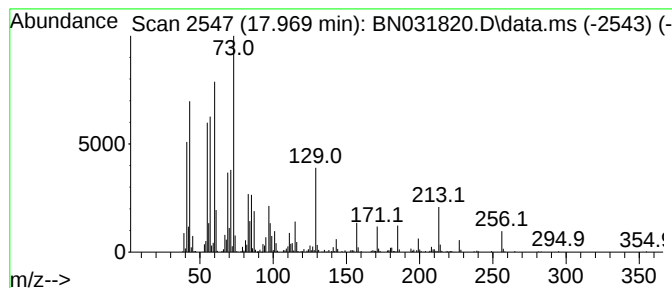
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TIC Integration Parameters: LSCINT.P

Peak Number 9 n-Hexadecanoic acid Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.969	4.86 ng/ul	1224940	Phenanthrene-d10	17.063

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		Tridecanoic acid	214	C13H26O2	000638-53-9	90



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ALS Vial : 9 Sample Multiplier: 1

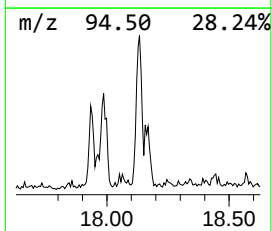
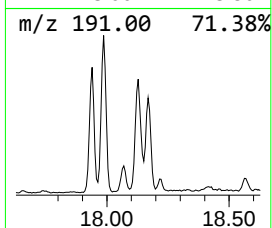
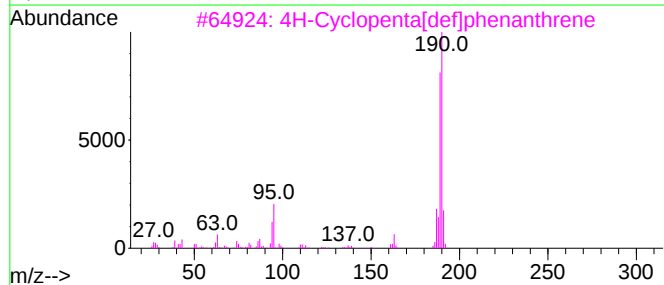
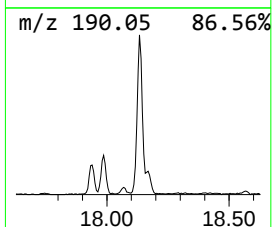
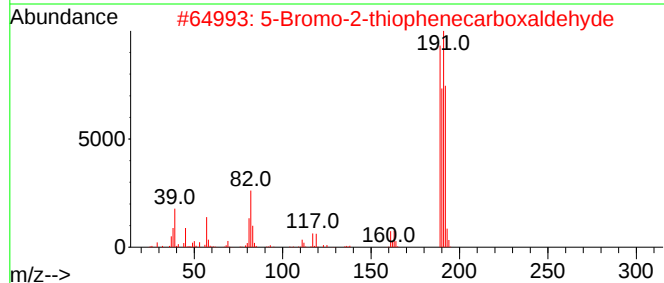
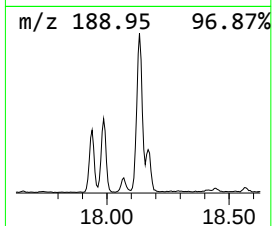
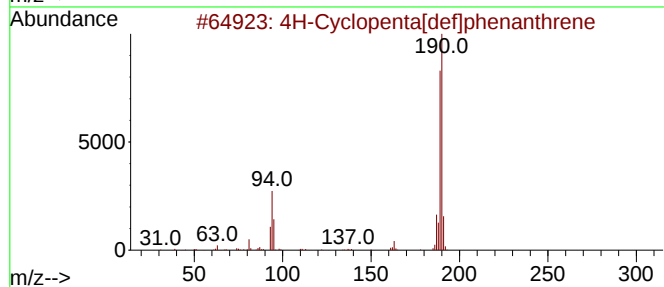
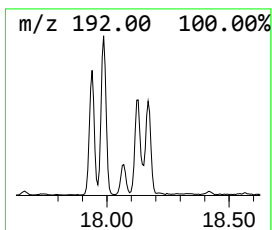
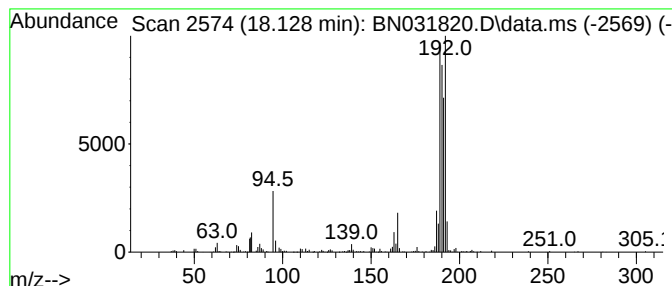
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ClientSampleId :
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 4H-Cyclopenta[def]phenanthrene Concentration Rank 21

R.T.	EstConc	Area	Relative to ISTD	R.T.	
18.128	2.08 ng/ul	523200	Phenanthrene-d10	17.063	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	60
2	5-Bromo-2-thiophenecarboxaldehyde	190	C5H3BrOS	004701-17-1	53
3	4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	49
4	6H-Cyclobuta[jk]phenanthrene	190	C15H10	083469-43-6	49
5	Naphtho[2,3-b]norbornadiene	192	C15H12	107426-38-0	42



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN060624\
Data File : BN031820.D
Acq On : 06 Jun 2024 15:55
Operator : MA/JU
Sample : P2634-06
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
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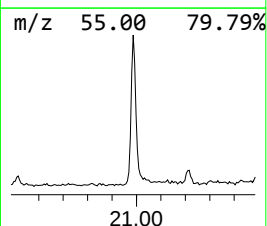
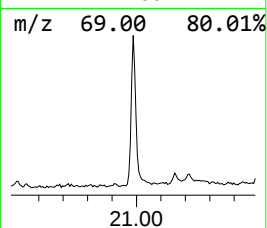
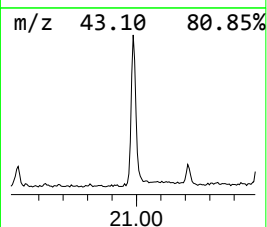
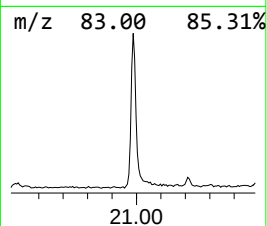
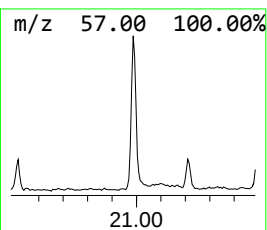
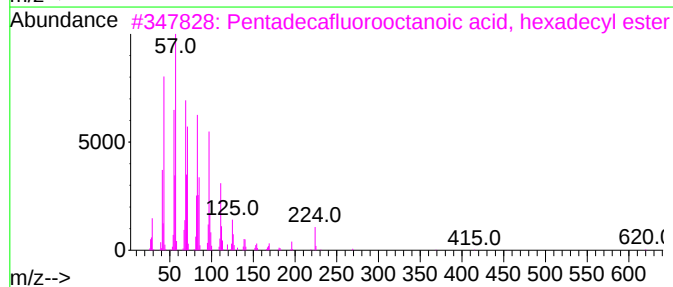
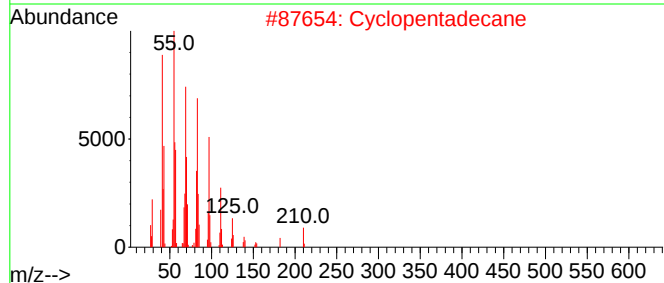
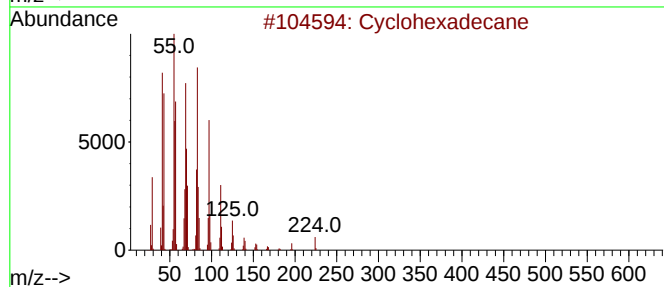
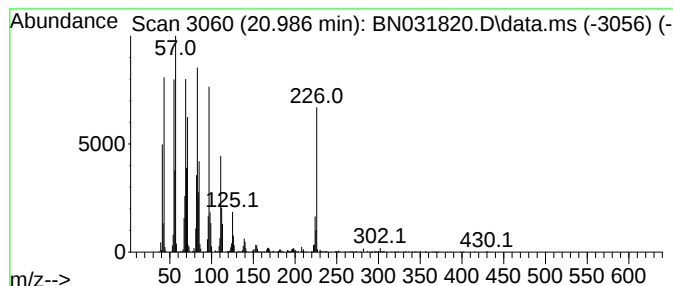
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Quant Title : SVOA CALIBRATION

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

Peak Number 11 (DEL) Alkane: Cyclic20.986 Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.986	3.92 ng/ul	1260160	Chrysene-d12	21.298

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexadecane	224	C16H32	000295-65-8	94
2			Cyclopentadecane	210	C15H30	000295-48-7	93
3			Pentadecafluorooctanoic acid, he...	638	C24H33F15O2	1000406-04-6	91
4			Heptafluorobutyric acid, pentade...	424	C19H31F7O2	959261-23-5	90
5			1-Heneicosanol	312	C21H44O	015594-90-8	89



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN060624\
Data File : BN031820.D
Acq On : 06 Jun 2024 15:55
Operator : MA/JU
Sample : P2634-06
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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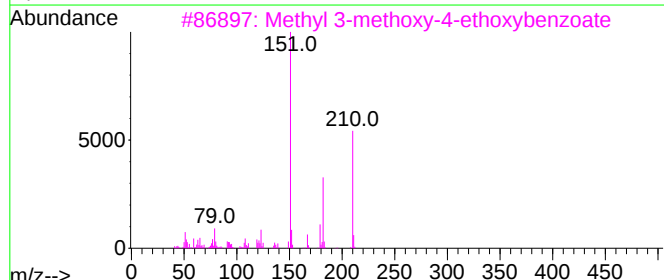
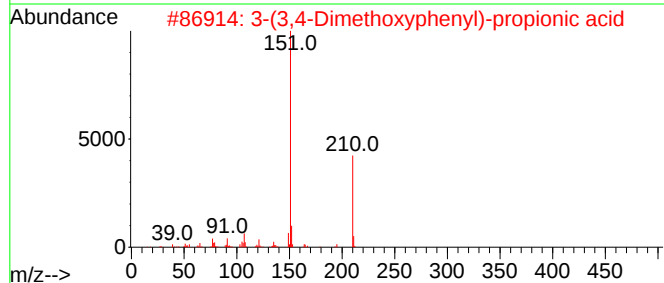
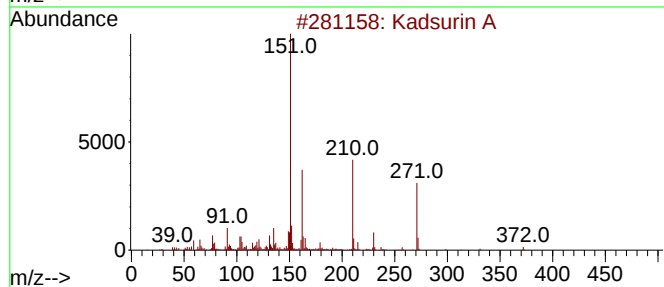
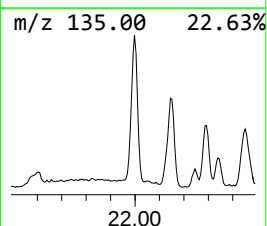
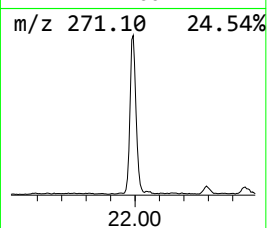
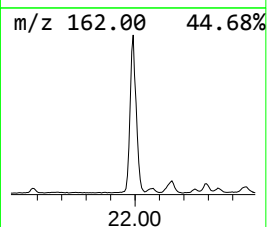
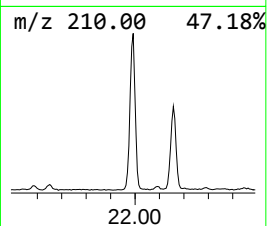
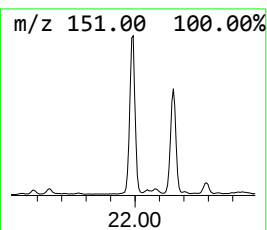
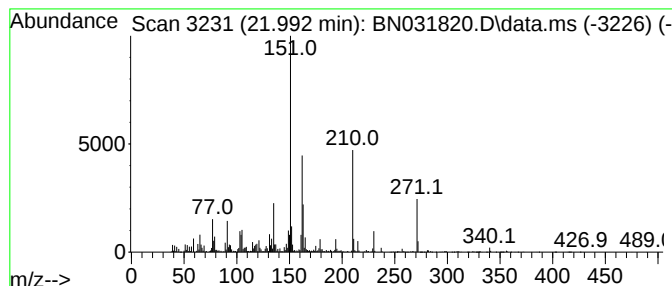
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Quant Title : SVOA CALIBRATION

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

Peak Number 12 Kadsurin A Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.992	11.58 ng/ul	3721670	Chrysene-d12	21.298

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Kadsurin A	372	C21H24O6	099340-07-5	76
2			3-(3,4-Dimethoxyphenyl)-propioni...	210	C11H14O4	002107-70-2	45
3			Methyl 3-methoxy-4-ethoxybenzoate	210	C11H14O4	003535-24-8	43
4			3-(3,4-Dimethoxyphenyl)-propioni...	210	C11H14O4	002107-70-2	43
5			Benzeneacetic acid, 3,4-dimethox...	210	C11H14O4	015964-79-1	43



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN060624\
Data File : BN031820.D
Acq On : 06 Jun 2024 15:55
Operator : MA/JU
Sample : P2634-06
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
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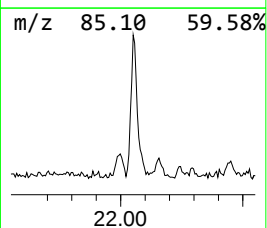
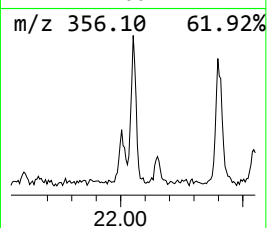
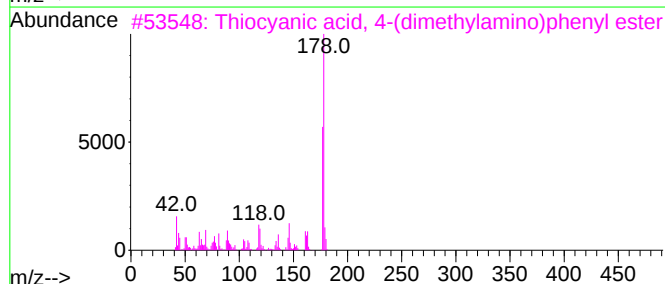
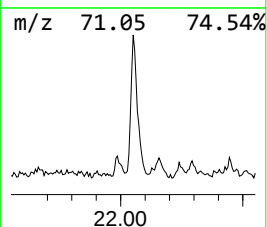
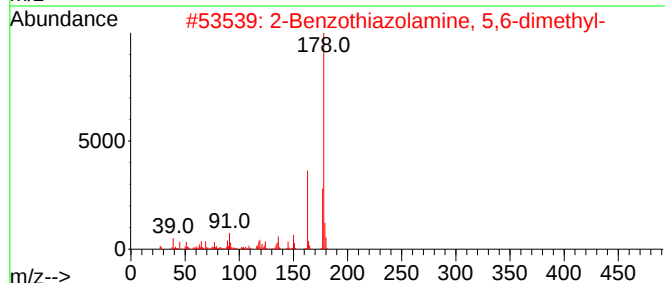
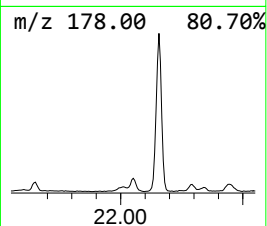
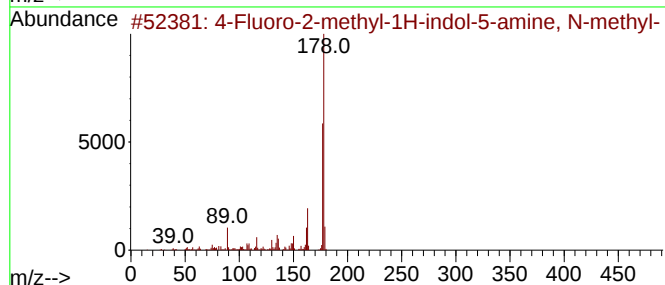
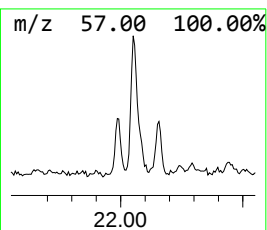
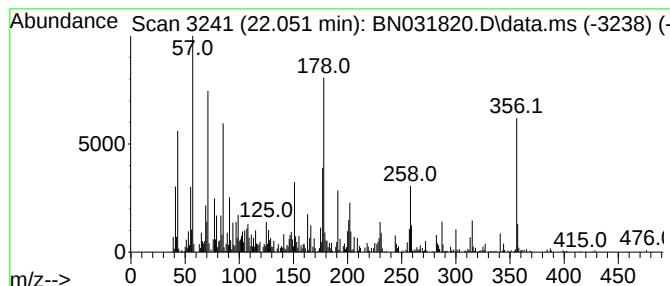
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Quant Title : SVOA CALIBRATION

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

Peak Number 13 unknown-01 Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.051	2.48 ng/ul	798009	Chrysene-d12	21.298

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			4-Fluoro-2-methyl-1H-indol-5-ami...	178	C10H11FN2	1010508-42-8	30
2			2-Benzothiazolamine, 5,6-dimethyl-	178	C9H10N2S	029927-08-0	25
3			Thiocyanic acid, 4-(dimethylamin...	178	C9H10N2S	007152-80-9	25
4			2-(4-Methoxyphenyl)-1,3,2-dioxab...	178	C9H11B03	069519-11-5	25
5			4-Aza-6-oxaocta-1,5-dione, 2-met...	311	C19H21N03	1000152-30-4	25



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN060624\
Data File : BN031820.D
Acq On : 06 Jun 2024 15:55
Operator : MA/JU
Sample : P2634-06
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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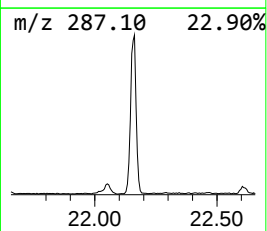
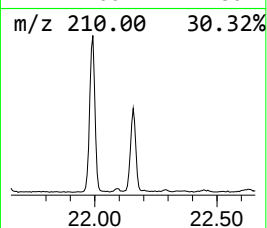
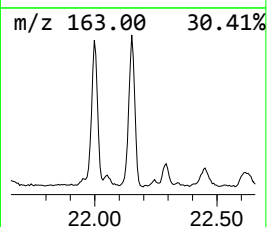
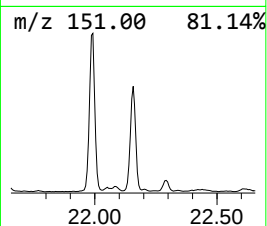
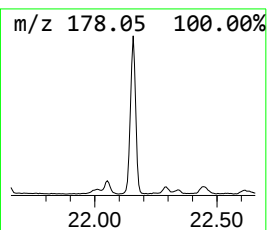
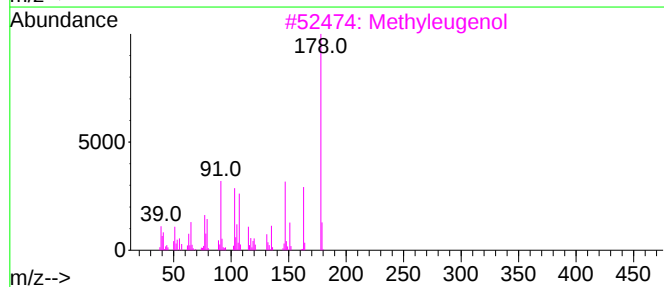
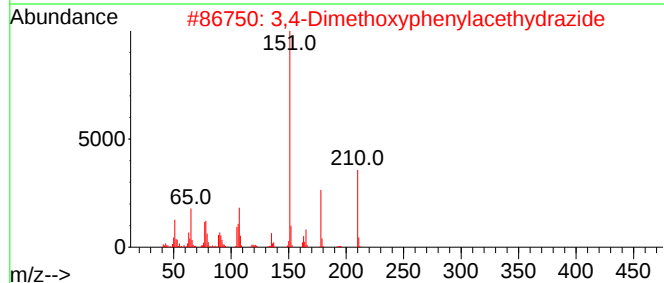
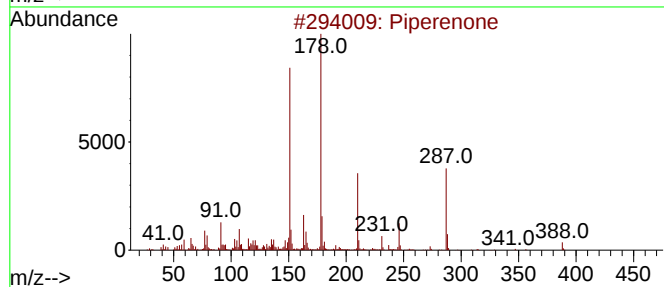
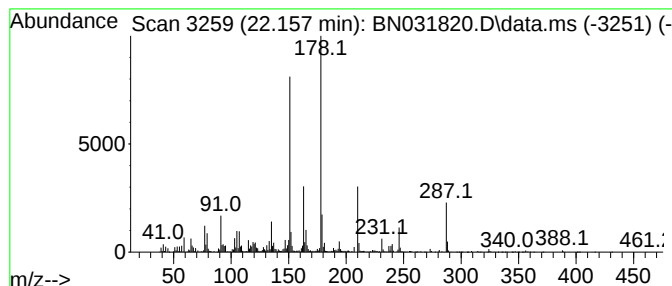
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Quant Title : SVOA CALIBRATION

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

Peak Number 14 Piperone Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.157	12.09 ng/ul	3883760	Chrysene-d12	21.298

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Piperone	388	C22H28O6	057625-31-7	90
2			3,4-Dimethoxyphenylacethydrazide	210	C10H14N2O3	060075-23-2	43
3			Methyleugenol	178	C11H14O2	000093-15-2	41
4			Methyleugenol	178	C11H14O2	000093-15-2	41
5			Methyleugenol	178	C11H14O2	000093-15-2	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN060624\
Data File : BN031820.D
Acq On : 06 Jun 2024 15:55
Operator : MA/JU
Sample : P2634-06
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
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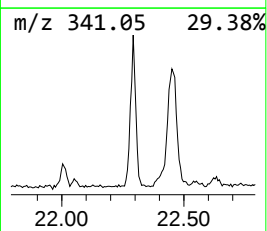
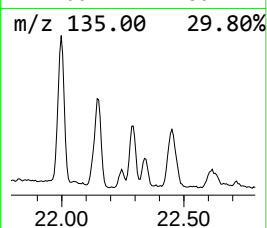
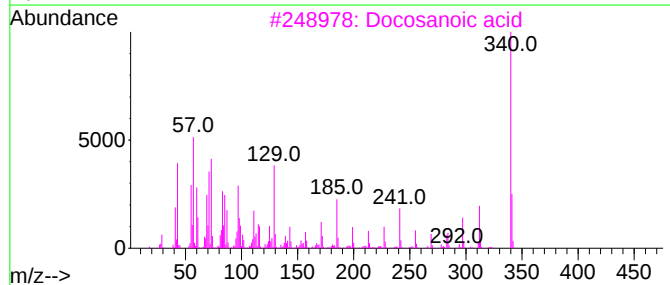
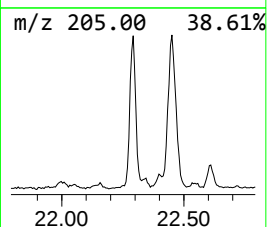
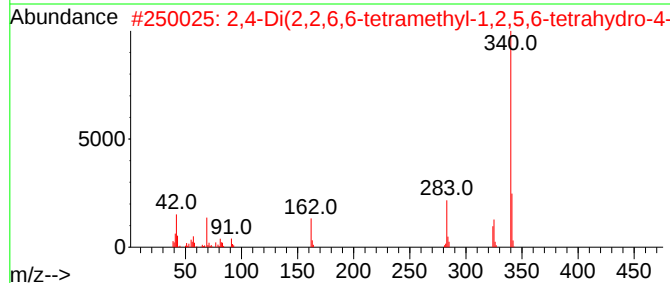
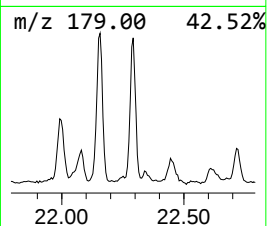
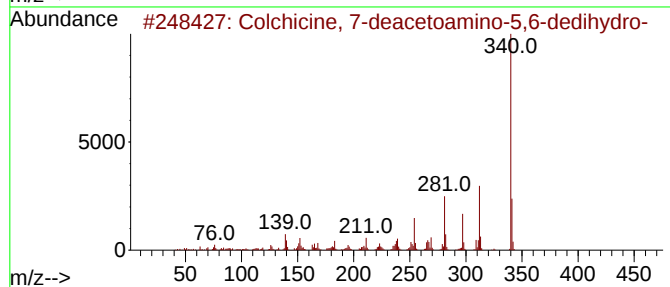
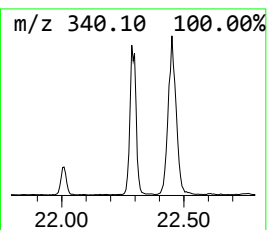
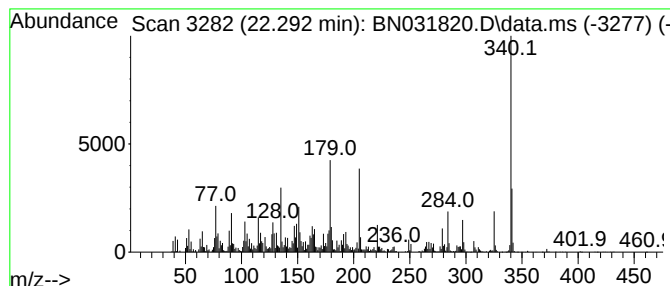
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Quant Title : SVOA CALIBRATION

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

Peak Number 15 unknown-02 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.292	6.62 ng/ul	2127290	Chrysene-d12	21.298

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Colchicine, 7-deacetoamino-5,6-d...	340	C20H20O5	076265-62-8	42
2			2,4-Di(2,2,6,6-tetramethyl-1,2,5...	341	C22H35N3	1000311-57-0	41
3			Docosanoic acid	340	C22H44O2	000112-85-6	41
4			Phenol, 2,6-bis(1,1-dimethylethy...	340	C23H32O2	020690-84-0	41
5			2-Phenyl-4,6-di(2-hydroxyphenyl)...	340	C22H16N2O2	052829-03-5	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN060624\
Data File : BN031820.D
Acq On : 06 Jun 2024 15:55
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Sample : P2634-06
Misc :
ALS Vial : 9 Sample Multiplier: 1

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ClientSampleId :
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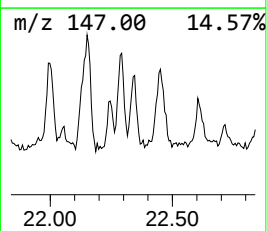
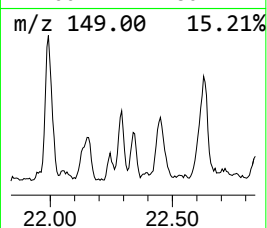
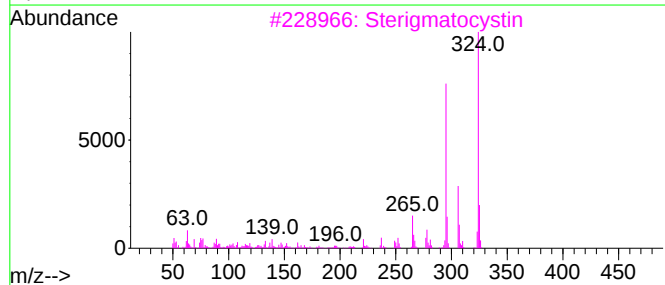
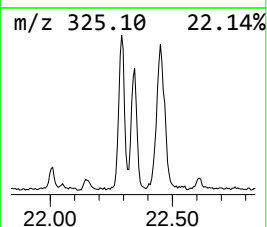
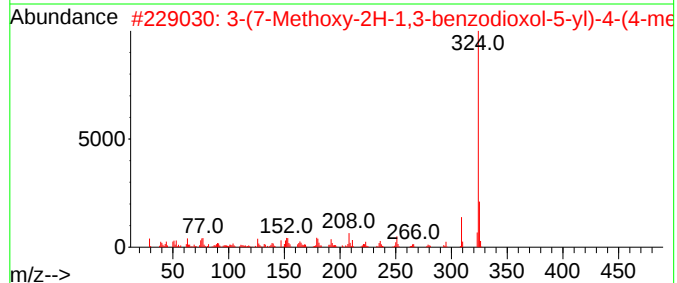
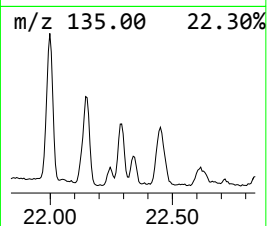
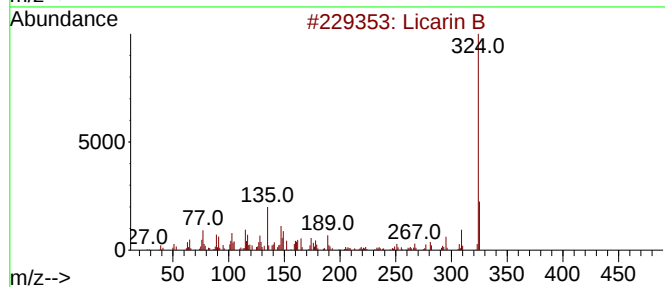
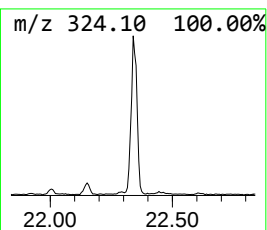
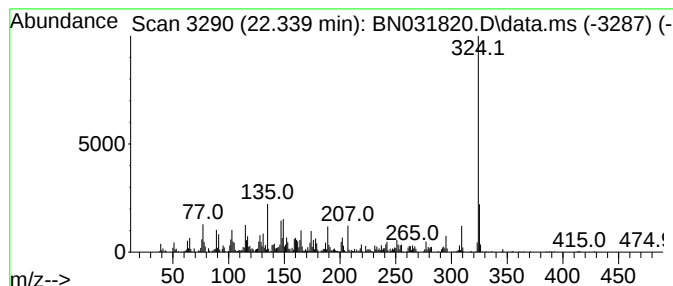
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Quant Title : SVOA CALIBRATION

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

Peak Number 16 Licarin B Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.339	2.43 ng/ul	781082	Chrysene-d12	21.298

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Licarin B	324	C20H20O4	051020-87-2	96
2			3-(7-Methoxy-2H-1,3-benzodioxol-...	324	C18H16N2O4	1000444-04-3	76
3			Sterigmatocystin	324	C18H12O6	010048-13-2	70
4			11H-Cyclopenta[a]phenanthrene-15...	324	C20H20O4	056588-57-9	64
5			4-(4-Hydroxyphenyl)-2-methyl-1(2...	324	C18H20N2O2Si	1000462-88-6	58



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Data File : BN031820.D
Acq On : 06 Jun 2024 15:55
Operator : MA/JU
Sample : P2634-06
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
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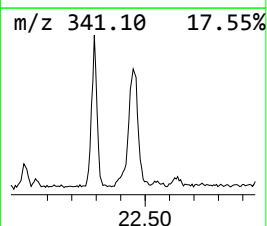
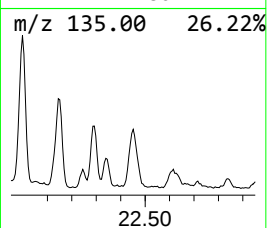
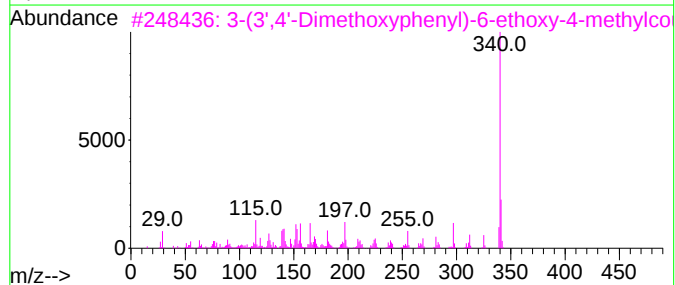
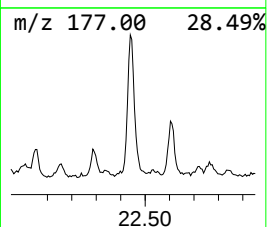
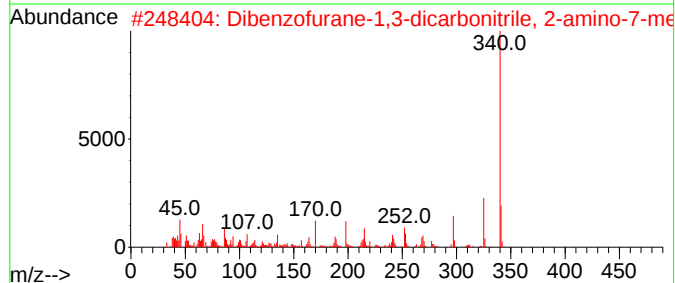
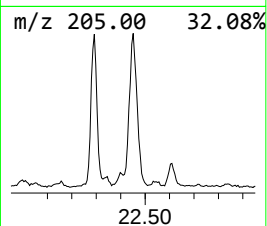
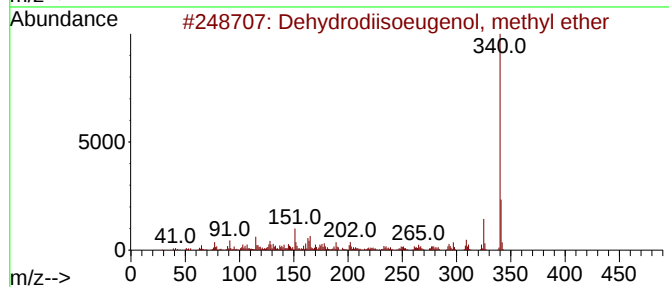
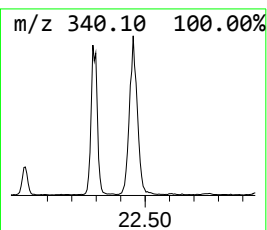
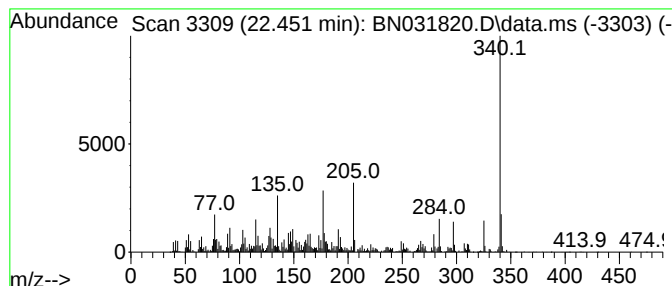
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Quant Title : SVOA CALIBRATION

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

Peak Number 17 Dehydrodiisoeugenol, methyl... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.451	8.91 ng/ul	2863350	Chrysene-d12	21.298

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Dehydrodiisoeugenol, methyl ether	340	C21H24O4	006544-71-4	55
2			Dibenzofurane-1,3-dicarbonitrile...	340	C20H12N4O2	330158-97-9	55
3			3-(3',4'-Dimethoxyphenyl)-6-etho...	340	C20H20O5	720674-30-6	49
4			Phenol, 2,2'-methylenebis[6-meth...	340	C21H24O4	055334-55-9	49
5			2,4-Di(2,2,6,6-tetramethyl-1,2,5...	341	C22H35N3	1000311-57-0	46



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN060624\
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Acq On : 06 Jun 2024 15:55
Operator : MA/JU
Sample : P2634-06
Misc :
ALS Vial : 9 Sample Multiplier: 1

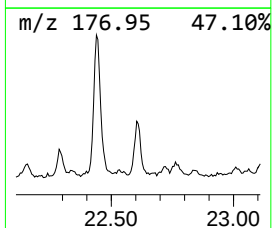
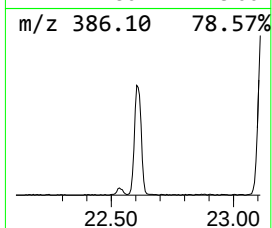
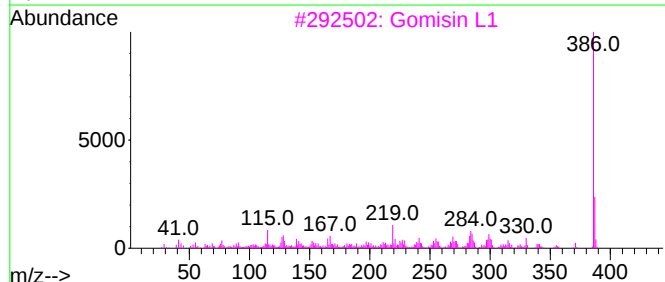
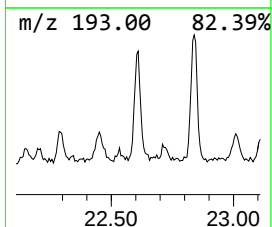
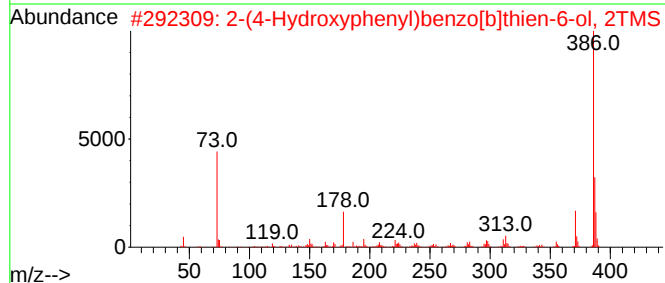
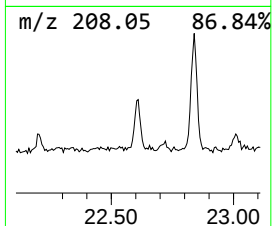
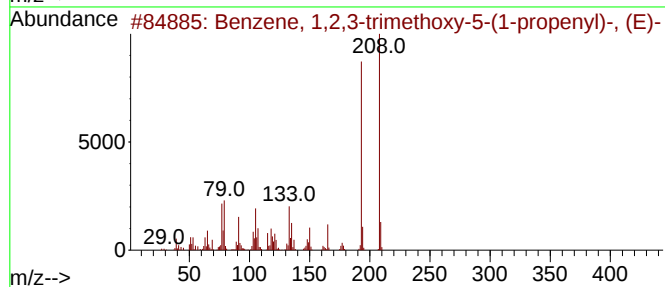
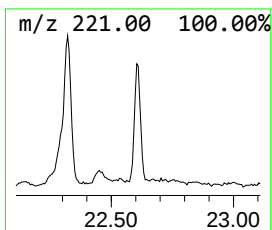
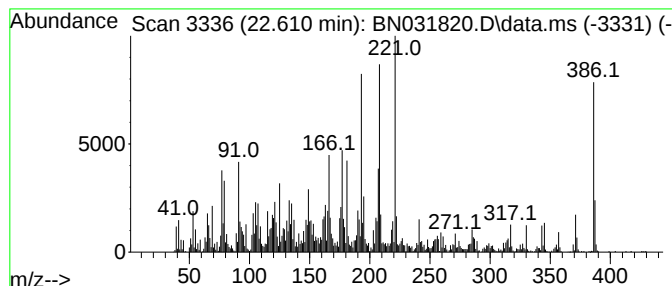
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Quant Title : SVOA CALIBRATION

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

Peak Number 18 Benzene, 1,2,3-trimethoxy-5... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.	
22.610	6.60 ng/ul	2119940	Chrysene-d12	21.298	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,2,3-trimethoxy-5-(1-p...	208	C12H16O3	005273-85-8	60
2	2-(4-Hydroxyphenyl)benzo[b]thien...	386	C20H26O2SSi2	1000499-61-8	55
3	Gomisin L1	386	C22H26O6	082425-43-2	53
4	Benzene, 1,2,3-trimethoxy-5-(2-p...	208	C12H16O3	000487-11-6	51
5	Benzene, 1,1'-(1,2-dimethyl-1,2-...	208	C16H16	000782-06-9	47



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN060624\
Data File : BN031820.D
Acq On : 06 Jun 2024 15:55
Operator : MA/JU
Sample : P2634-06
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
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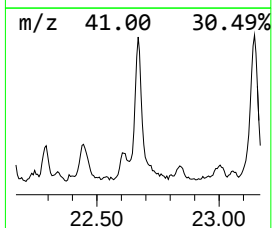
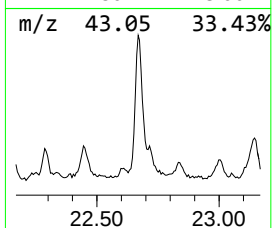
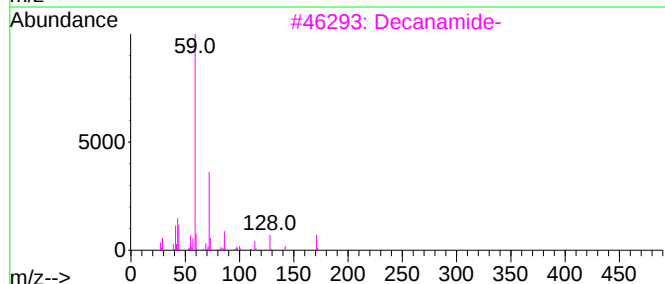
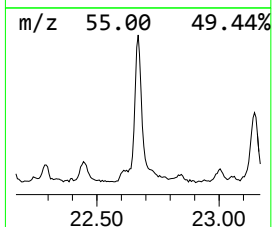
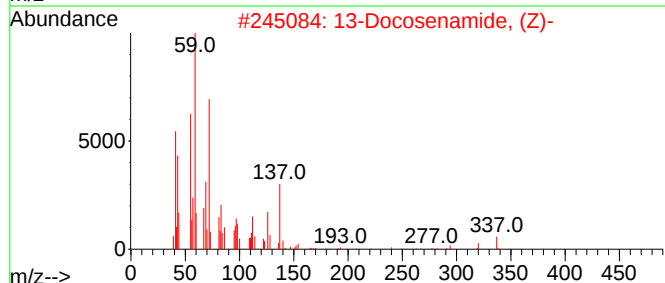
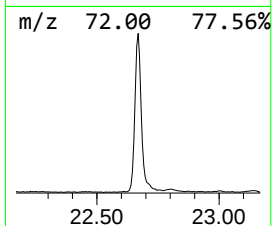
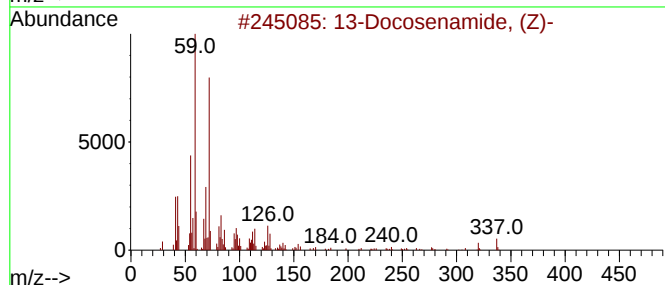
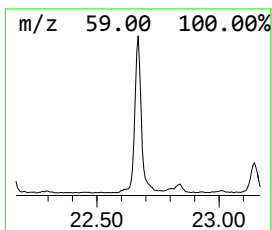
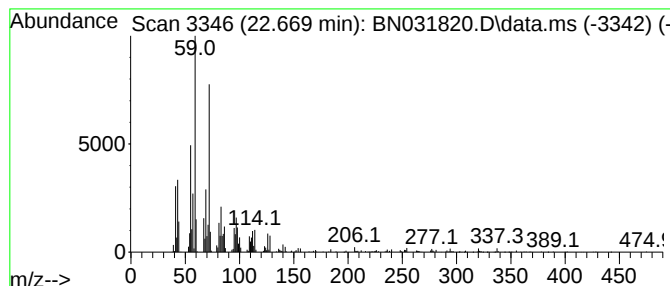
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Quant Title : SVOA CALIBRATION

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

Peak Number 19 13-Docosenamide, (Z)- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.669	3.50 ng/ul	1124070	Chrysene-d12	21.298

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			13-Docosenamide, (Z)-	337	C22H43NO	000112-84-5	97
2			13-Docosenamide, (Z)-	337	C22H43NO	000112-84-5	72
3			Decanamide-	171	C10H21NO	002319-29-1	53
4			9-Octadecenamide	281	C18H35NO	003322-62-1	47
5			Hexadecanamide	255	C16H33NO	000629-54-9	47



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN060624\
Data File : BN031820.D
Acq On : 06 Jun 2024 15:55
Operator : MA/JU
Sample : P2634-06
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
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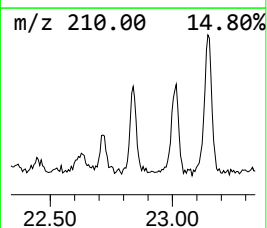
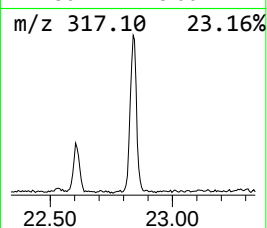
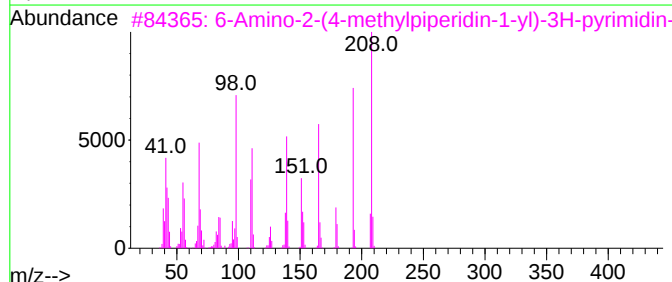
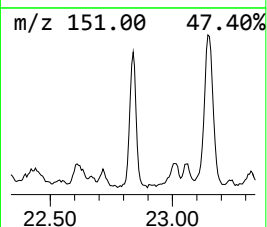
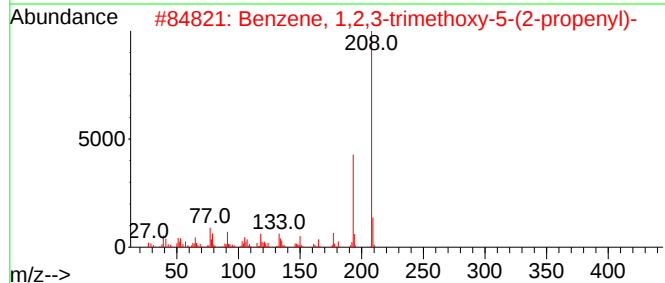
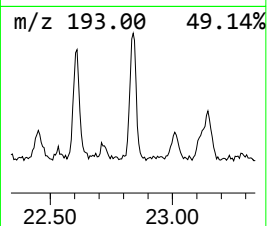
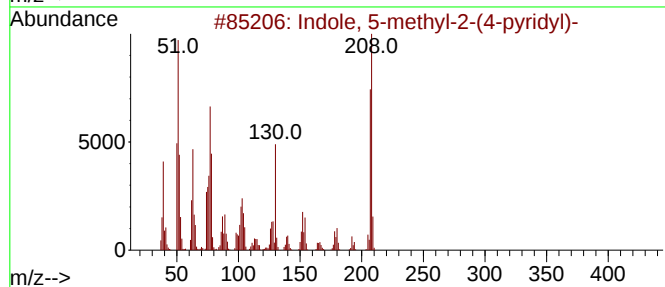
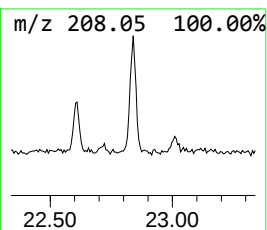
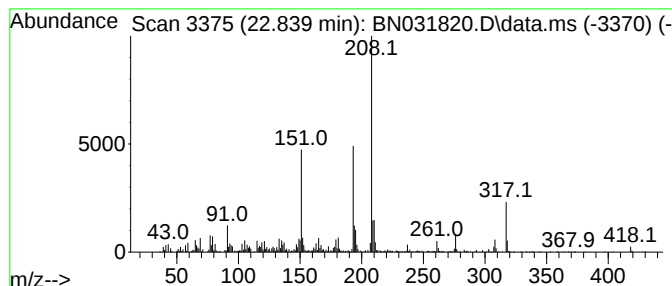
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Quant Title : SVOA CALIBRATION

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

Peak Number 20 Indole, 5-methyl-2-(4-pyrid... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.839	5.94 ng/ul	1080940	Perylene-d12	24.239

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Indole, 5-methyl-2-(4-pyridyl)-	208	C14H12N2	107919-92-6	70
2			Benzene, 1,2,3-trimethoxy-5-(2-p...	208	C12H16O3	000487-11-6	60
3			6-Amino-2-(4-methylpiperidin-1-y...	208	C10H16N4O	1000388-65-9	60
4			Benzene, 1,2,3-trimethoxy-5-(2-p...	208	C12H16O3	000487-11-6	55
5			Asarone	208	C12H16O3	002883-98-9	55



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN060624\
Data File : BN031820.D
Acq On : 06 Jun 2024 15:55
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Sample : P2634-06
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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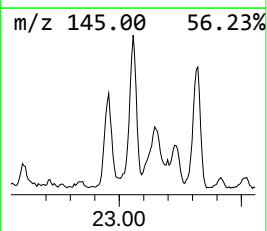
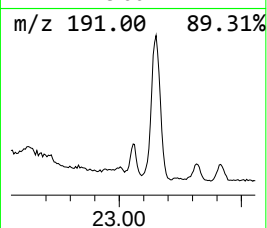
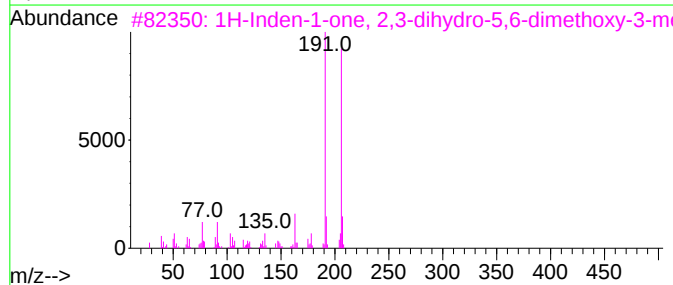
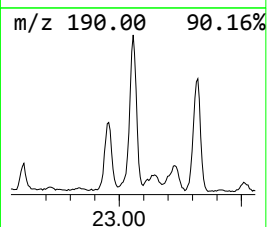
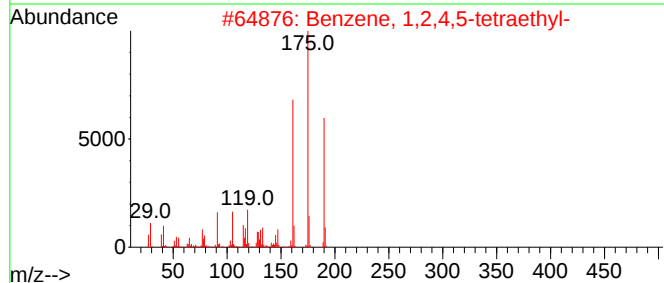
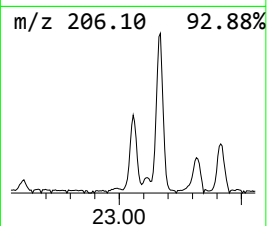
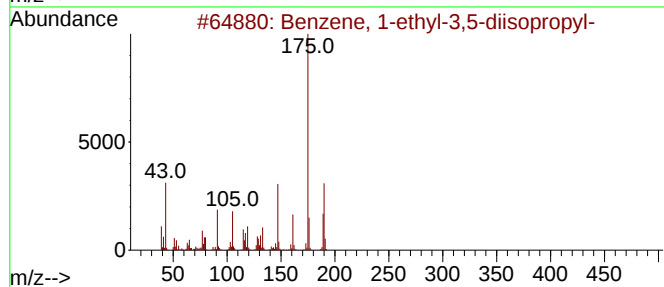
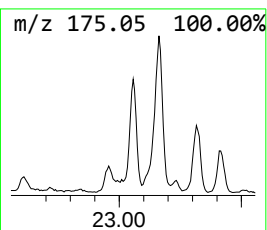
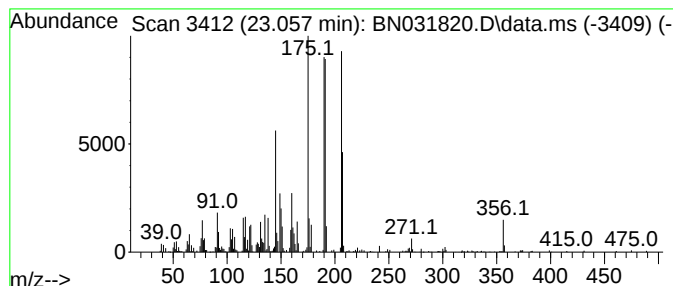
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Quant Title : SVOA CALIBRATION

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

Peak Number 21 unknown-03 Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.057	3.17 ng/ul	576601	Perylene-d12	24.239

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethyl-3,5-diisopropyl-	190	C14H22	015181-13-2	38
2			Benzene, 1,2,4,5-tetraethyl-	190	C14H22	000635-81-4	38
3			1H-Inden-1-one, 2,3-dihydro-5,6-...	206	C12H14O3	004082-25-1	38
4			3,4-Dihydroisoquinolin-6-ol, 7-m...	191	C11H13NO2	004602-70-4	30
5			1-Ethyl-5-fluoro-1H-indole-3-car...	191	C11H10FNO	620175-69-1	30



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN060624\
Data File : BN031820.D
Acq On : 06 Jun 2024 15:55
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Sample : P2634-06
Misc :
ALS Vial : 9 Sample Multiplier: 1

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ClientSampleId :
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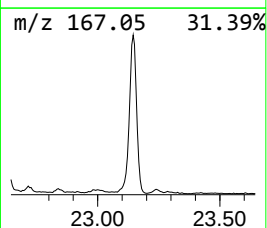
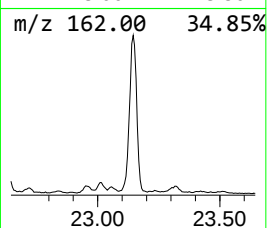
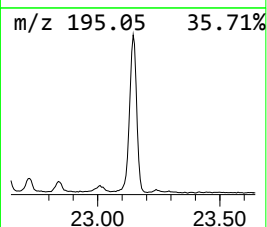
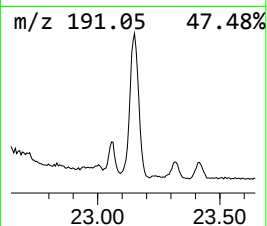
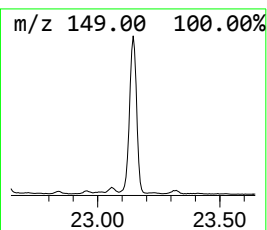
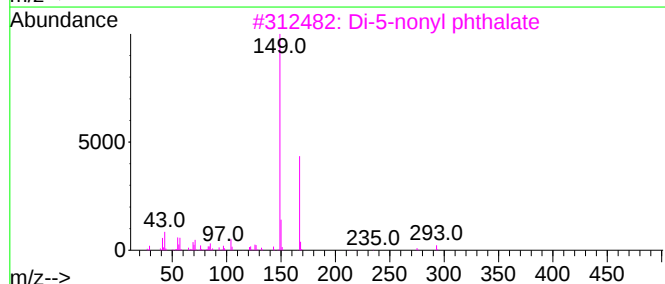
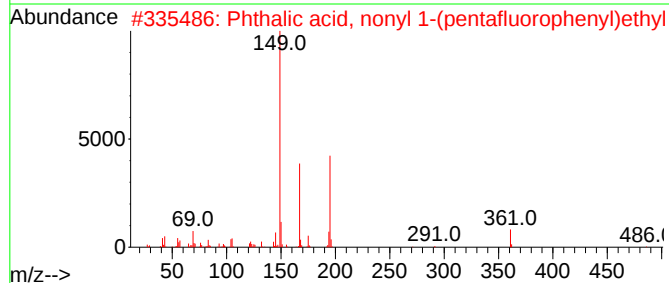
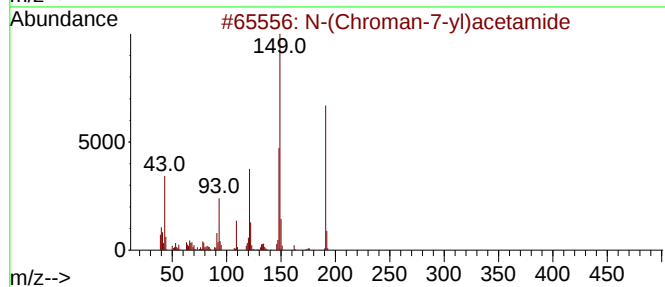
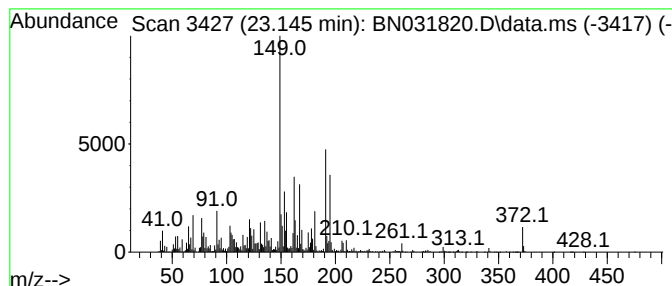
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Quant Title : SVOA CALIBRATION

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

Peak Number 22 unknown-04 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.145	49.98 ng/ul	9098180	Perylene-d12	24.239

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			N-(Chroman-7-yl)acetamide	191	C11H13NO2	1000306-69-7	46
2			Phthalic acid, nonyl 1-(pentaflu...	486	C25H27F5O4	1000382-93-2	38
3			Di-5-nonyl phthalate	418	C26H42O4	094054-23-6	35
4			Bis(2-ethylhexyl) phthalate	390	C24H38O4	000117-81-7	35
5			1H-[1,2,3]Triazole-4-carboxylic ...	372	C15H16N8O4	1000303-81-5	30



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN060624\
Data File : BN031820.D
Acq On : 06 Jun 2024 15:55
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Sample : P2634-06
Misc :
ALS Vial : 9 Sample Multiplier: 1

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Propylene Glycol	3.446	2.7	ng/ul	314747	1	7.669	2319730	20.0
.alpha.-Pinene	6.358	8.9	ng/ul	1029260	1	7.669	2319730	20.0
Camphene	6.663	3.3	ng/ul	384511	1	7.669	2319730	20.0
.beta.-Pinene	7.105	15.2	ng/ul	1757060	1	7.669	2319730	20.0
1,3,6-Octatrien...	7.563	2.0	ng/ul	235276	1	7.669	2319730	20.0
D-Limonene	7.887	2.1	ng/ul	247282	1	7.669	2319730	20.0
1-Phenoxypropan...	11.199	2.4	ng/ul	408580	2	10.451	3370940	20.0
.gamma.-Asarone	14.869	2.6	ng/ul	552678	3	14.316	4297300	20.0
n-Hexadecanoic ...	17.969	4.9	ng/ul	1224940	4	17.063	5041660	20.0
4H-Cyclopenta[d...	18.128	2.1	ng/ul	523200	4	17.063	5041660	20.0
(DEL) Alkane: C...	20.986	3.9	ng/ul	1260160	5	21.298	6427310	20.0
Kadsurin A	21.992	11.6	ng/ul	3721670	5	21.298	6427310	20.0
unknown-01	22.051	2.5	ng/ul	798009	5	21.298	6427310	20.0
Piperenone	22.157	12.1	ng/ul	3883760	5	21.298	6427310	20.0
unknown-02	22.292	6.6	ng/ul	2127290	5	21.298	6427310	20.0
Licarin B	22.339	2.4	ng/ul	781082	5	21.298	6427310	20.0
Dehydrodiisoeug...	22.451	8.9	ng/ul	2863350	5	21.298	6427310	20.0
Benzene, 1,2,3-...	22.610	6.6	ng/ul	2119940	5	21.298	6427310	20.0
13-Docosenamide...	22.669	3.5	ng/ul	1124070	5	21.298	6427310	20.0
Indole, 5-methy...	22.839	5.9	ng/ul	1080940	6	24.239	3641040	20.0
unknown-03	23.057	3.2	ng/ul	576601	6	24.239	3641040	20.0
unknown-04	23.145	50.0	ng/ul	9098180	6	24.239	3641040	20.0