

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN060624\
 Data File : BN031821.D
 Acq On : 06 Jun 2024 16:34
 Operator : MA/JU
 Sample : P2634-07
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 BHBX1

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: BN031821.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.187	29	34	42	rVB	80150	113412	1.36%	0.121%
2	3.446	74	78	88	rBV	187923	266649	3.19%	0.284%
3	3.587	99	102	116	rBV	998181	1457787	17.44%	1.550%
4	3.905	151	156	161	rBV	19994	29473	0.35%	0.031%
5	4.805	303	309	317	rVB2	10175	21010	0.25%	0.022%
6	5.769	468	473	483	rBV	55687	94760	1.13%	0.101%
7	6.358	568	573	580	rVB	107187	170206	2.04%	0.181%
8	6.664	619	625	631	rVB2	47137	81534	0.98%	0.087%
9	6.846	644	656	667	rBV	1389639	2297886	27.49%	2.443%
10	7.011	678	684	694	rVB	1139724	1788855	21.40%	1.902%
11	7.105	694	700	710	rVB	147332	251960	3.01%	0.268%
12	7.205	710	717	726	rBV	1412522	2235830	26.75%	2.377%
13	7.281	726	730	736	rVV2	35503	55245	0.66%	0.059%
14	7.563	772	778	783	rBV	29341	47550	0.57%	0.051%
15	7.669	788	796	803	rBV	1336253	2141139	25.61%	2.277%
16	7.740	803	808	815	rVV2	50325	91600	1.10%	0.097%
17	7.799	815	818	825	rVB4	7942	15197	0.18%	0.016%
18	7.887	827	833	840	rVV3	26357	41060	0.49%	0.044%
19	8.046	853	860	866	rBV10	5773	12460	0.15%	0.013%
20	8.375	909	916	925	rBV	1805480	2873111	34.37%	3.055%
21	8.828	984	993	1001	rBV	1303436	2139527	25.60%	2.275%
22	8.987	1016	1020	1026	rVB7	8461	16709	0.20%	0.018%
23	9.228	1052	1061	1067	rBV2	19605	31793	0.38%	0.034%
24	9.393	1083	1089	1100	rBV	176031	278911	3.34%	0.297%
25	9.546	1103	1115	1127	rBV	1031666	1767831	21.15%	1.880%
26	9.775	1149	1154	1161	rBV9	9591	20043	0.24%	0.021%
27	10.075	1197	1205	1216	rBV	1644282	2817182	33.70%	2.996%
28	10.452	1260	1269	1281	rVB	1845660	3148038	37.66%	3.347%
29	10.593	1283	1293	1309	rBV	1613726	2824622	33.79%	3.003%
30	10.834	1331	1334	1341	rVB7	6474	11042	0.13%	0.012%
31	11.004	1356	1363	1373	rVB	55029	106817	1.28%	0.114%
32	11.199	1390	1396	1412	rBV	175748	359359	4.30%	0.382%
33	11.822	1494	1502	1507	rBV4	10961	20166	0.24%	0.021%
34	11.881	1507	1512	1517	rVV2	7443	10946	0.13%	0.012%
35	12.046	1532	1540	1546	rBV2	32121	57051	0.68%	0.061%
36	12.657	1637	1644	1649	rBV4	11041	26106	0.31%	0.028%

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Integrator: RTE

Smoothing : OFF

Filtering: 5

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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
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37	13.040	1704	1709	1712	rBV6	10137	19881	0.24%	0.021%
38	13.140	1722	1726	1729	rBV2	17074	26348	0.32%	0.028%
39	13.181	1731	1733	1739	rVV6	13071	18338	0.22%	0.019%
40	13.275	1747	1749	1754	rVB4	8777	13304	0.16%	0.014%

41	13.504	1780	1788	1791	rBV7	5777	14208	0.17%	0.015%
42	13.634	1806	1810	1814	rVV6	15111	24442	0.29%	0.026%
43	13.728	1820	1826	1836	rBV	2733099	3920972	46.91%	4.169%
44	13.804	1836	1839	1841	rVV4	14238	19650	0.24%	0.021%
45	13.845	1842	1846	1854	rVB3	44726	69219	0.83%	0.074%

46	14.010	1862	1874	1886	rBV	3288642	5112657	61.16%	5.436%
47	14.157	1895	1899	1905	rBV9	5372	10774	0.13%	0.011%
48	14.222	1905	1910	1913	rBV4	11053	16823	0.20%	0.018%
49	14.316	1919	1926	1933	rVB	2704055	4040645	48.34%	4.297%
50	14.387	1933	1938	1949	rVB9	24542	67695	0.81%	0.072%

51	14.522	1953	1961	1971	rBV	1345677	2168519	25.94%	2.306%
52	14.745	1995	1999	2004	rVB5	19837	32429	0.39%	0.034%
53	14.869	2015	2020	2024	rVB	54466	76354	0.91%	0.081%
54	15.110	2058	2061	2068	rVB6	15280	22022	0.26%	0.023%
55	15.175	2068	2072	2075	rBV2	12584	17367	0.21%	0.018%

56	15.310	2088	2095	2102	rBV	4080890	6021844	72.04%	6.403%
57	15.363	2102	2104	2111	rVV2	15553	26180	0.31%	0.028%
58	15.440	2111	2117	2130	rVV	1053757	1546149	18.50%	1.644%
59	15.551	2132	2136	2145	rVB3	23713	49466	0.59%	0.053%
60	15.657	2151	2154	2161	rVB8	7262	16194	0.19%	0.017%

61	15.792	2172	2177	2182	rBV7	17007	29747	0.36%	0.032%
62	15.881	2187	2192	2197	rBV8	10687	23586	0.28%	0.025%
63	15.939	2197	2202	2206	rBV7	14768	21837	0.26%	0.023%
64	16.039	2214	2219	2226	rBV5	21815	41834	0.50%	0.044%
65	16.381	2268	2277	2279	rBV8	8546	23049	0.28%	0.025%

66	16.475	2289	2293	2295	rBV5	8388	12115	0.14%	0.013%
67	16.616	2311	2317	2321	rBV7	14741	34296	0.41%	0.036%
68	16.663	2323	2325	2330	rVB4	9620	13308	0.16%	0.014%
69	16.728	2331	2336	2342	rVB5	15870	31672	0.38%	0.034%
70	16.787	2342	2346	2348	rBV4	9035	13241	0.16%	0.014%

71	16.986	2376	2380	2384	rBV5	13997	22275	0.27%	0.024%
72	17.063	2386	2393	2398	rVV	3350171	4827720	57.75%	5.133%
73	17.104	2398	2400	2404	rVV	262368	311723	3.73%	0.331%
74	17.163	2404	2410	2420	rVV	4313847	6137909	73.43%	6.527%
75	17.263	2421	2427	2431	rVV5	28893	45584	0.55%	0.048%

76	17.445	2455	2458	2459	rBV3	12323	11988	0.14%	0.013%
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Integration Parameters: LSCINT.P

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

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

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

77	17.469	2460	2462	2466	rVB	18628	22607	0.27%	0.024%
78	17.645	2489	2492	2495	rBV4	11623	15569	0.19%	0.017%
79	17.739	2506	2508	2512	rVB5	14584	17012	0.20%	0.018%
80	17.963	2538	2546	2561	rBV2	449227	983916	11.77%	1.046%
81	18.128	2570	2574	2579	rBV2	63437	110428	1.32%	0.117%
82	18.169	2579	2581	2586	rVB	39055	49599	0.59%	0.053%
83	18.292	2598	2602	2606	rBV3	44848	69207	0.83%	0.074%
84	18.445	2624	2628	2632	rBV2	51372	75620	0.90%	0.080%
85	18.486	2632	2635	2640	rVB2	33438	46102	0.55%	0.049%
86	18.863	2695	2699	2704	rBV2	68648	102845	1.23%	0.109%
87	19.022	2719	2726	2733	rBV3	75933	149675	1.79%	0.159%
88	19.128	2735	2744	2752	rVV	522281	851909	10.19%	0.906%
89	19.251	2761	2765	2773	rBV2	158131	258323	3.09%	0.275%
90	19.463	2795	2801	2815	rBV	5637497	8359057	100.00%	8.888%
91	19.986	2888	2890	2893	rVB2	72449	64797	0.78%	0.069%
92	20.986	3056	3060	3067	rBV	744094	1019281	12.19%	1.084%
93	21.298	3106	3113	3118	rBV2	3316820	5151378	61.63%	5.478%
94	21.986	3226	3230	3236	rVB3	495403	878897	10.51%	0.935%
95	22.151	3255	3258	3265	rVB3	441708	662617	7.93%	0.705%
96	22.286	3277	3281	3286	rVB2	244337	373470	4.47%	0.397%
97	22.663	3341	3345	3352	rBV3	491854	862449	10.32%	0.917%
98	23.139	3421	3426	3435	rVB4	1022892	2166516	25.92%	2.304%
99	24.033	3570	3578	3587	rBV	2173388	5123182	61.29%	5.448%
100	24.233	3604	3612	3622	rBV	1524639	3953953	47.30%	4.204%

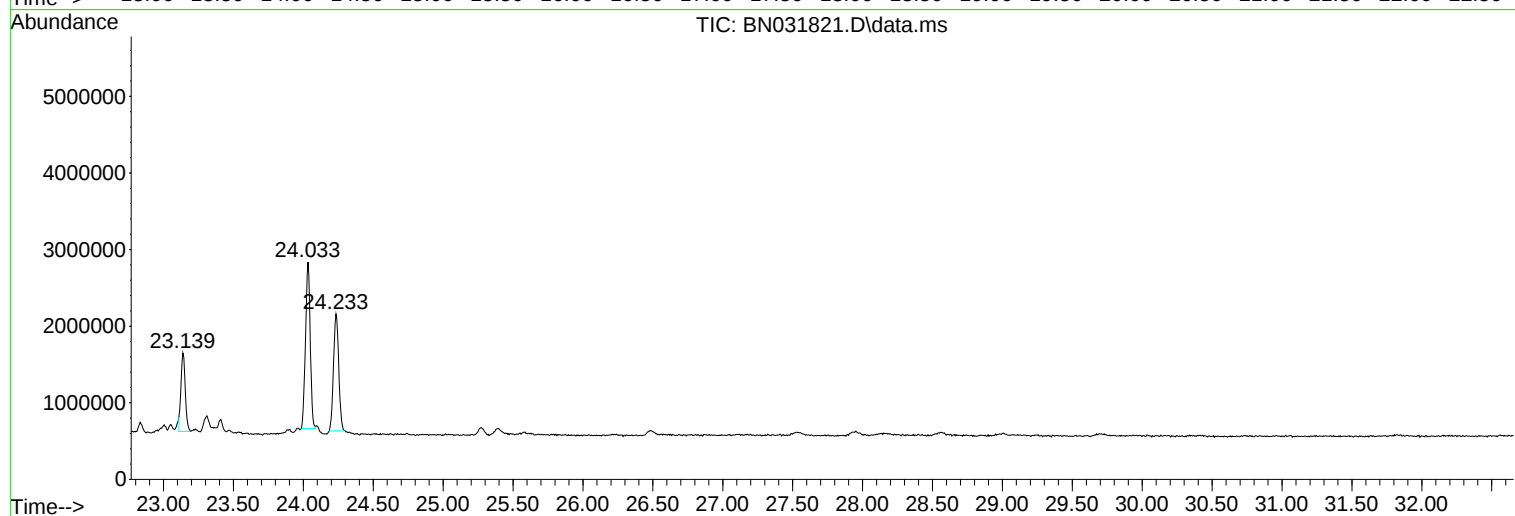
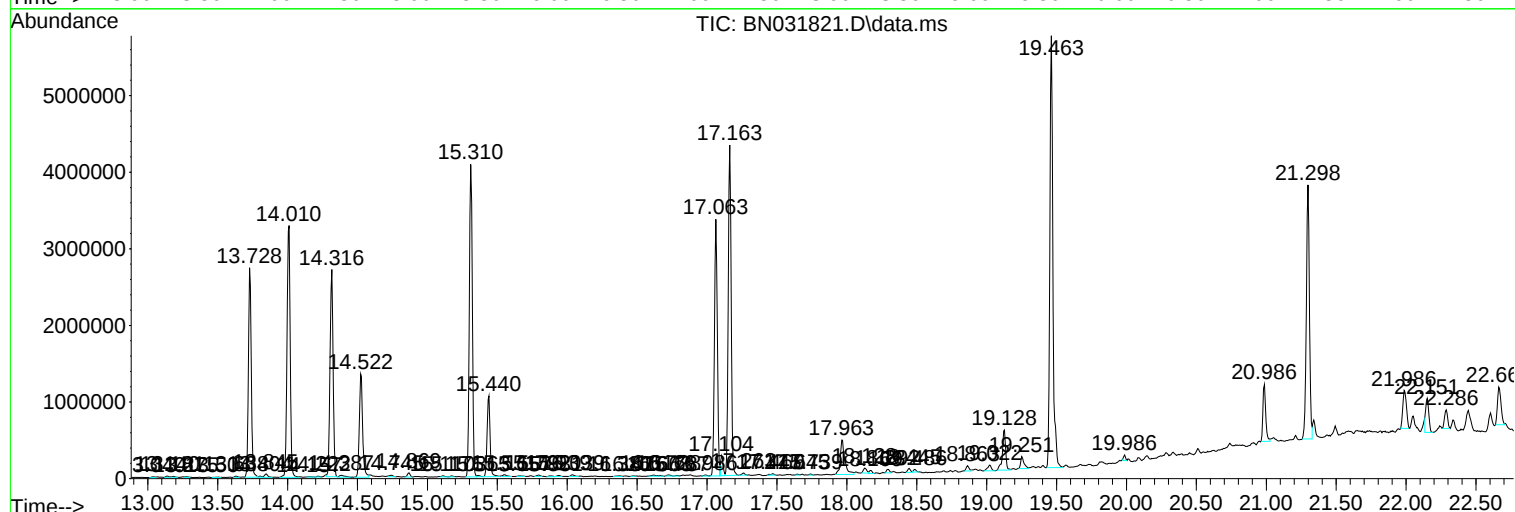
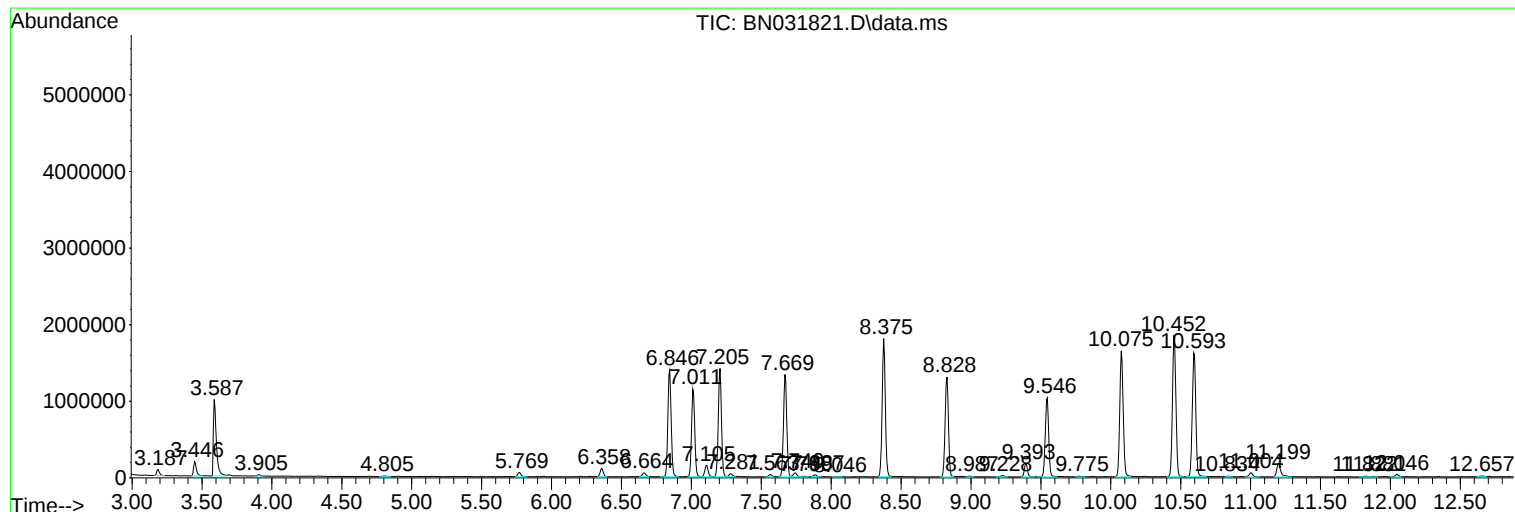
Sum of corrected areas: 94044640

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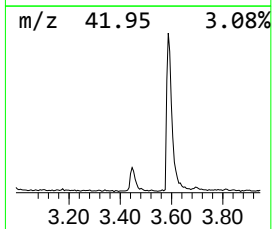
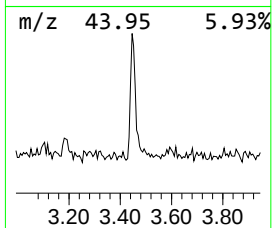
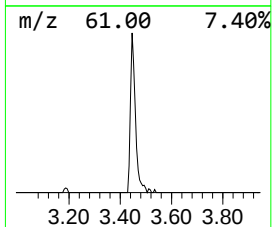
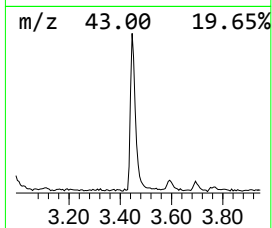
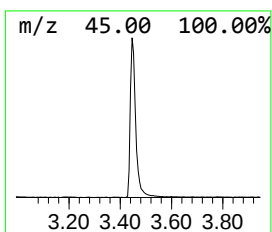
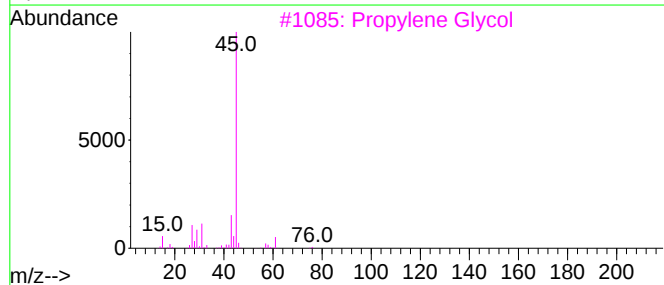
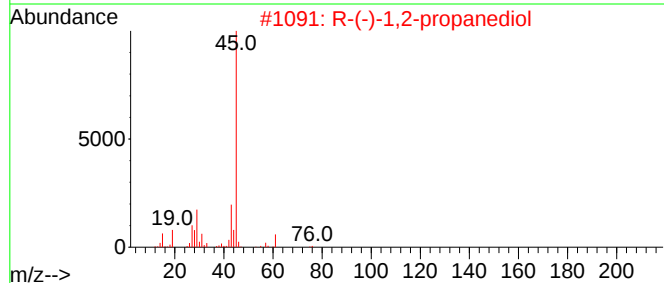
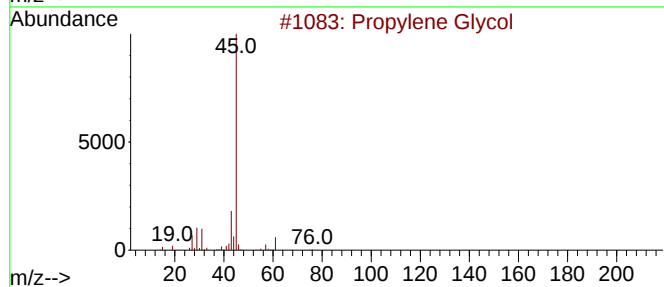
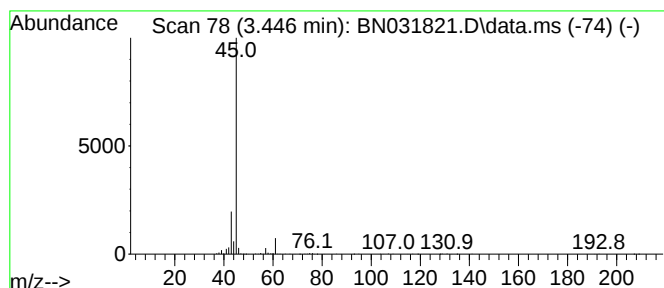
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN052924.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Propylene Glycol Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.446	2.49 ng/ul	266649	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			Propylene Glycol	76	C3H8O2	000057-55-6	56
4			Propanoic acid, 2-hydroxy-, meth...	104	C4H8O3	000547-64-8	40
5			(S)-(+)-1,2-Propanediol	76	C3H8O2	004254-15-3	40



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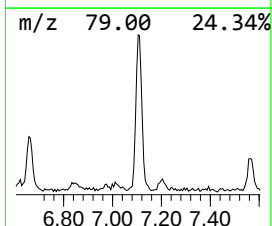
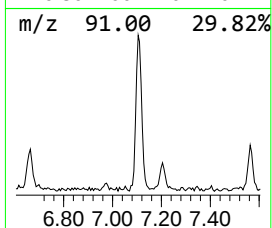
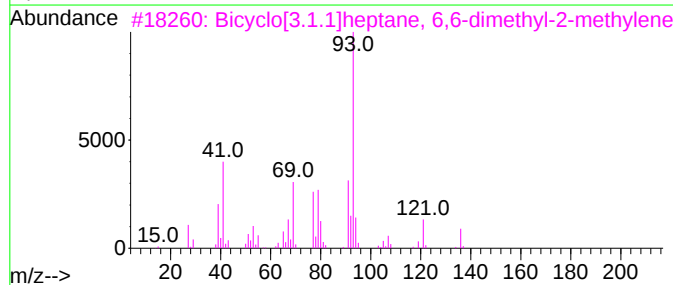
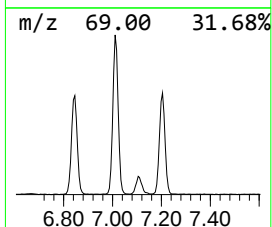
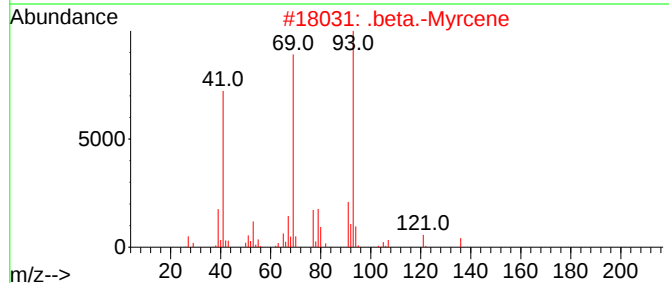
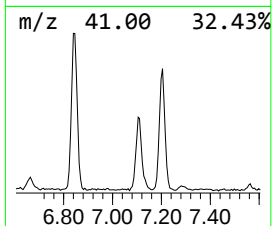
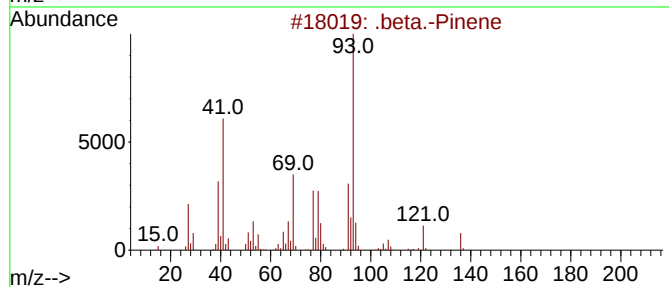
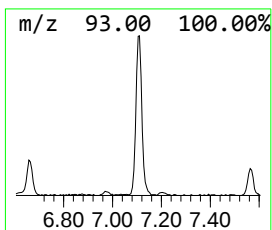
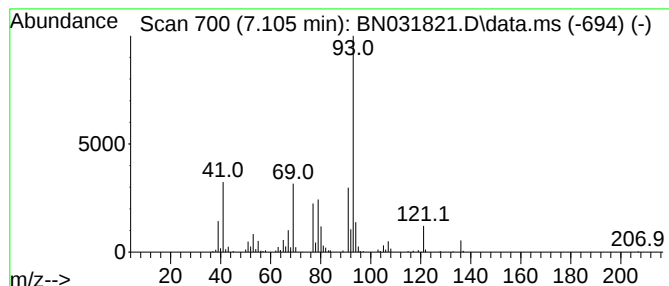
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN052924.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 .beta.-Pinene Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.105	2.35 ng/ul	251960	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	.	beta.-Myrcene	136	C10H16	000123-35-3	91	
3	Bicyclo[3.1.1]heptane, 6,6-dimet...	136	C10H16	018172-67-3	91		
4	.	beta.-Pinene	136	C10H16	000127-91-3	91	
5	Bicyclo[3.1.1]heptane, 6,6-dimet...	136	C10H16	018172-67-3	91		



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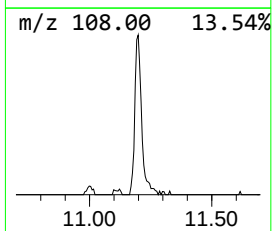
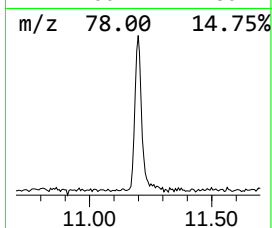
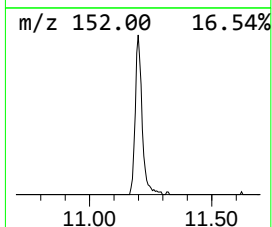
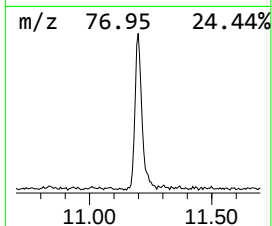
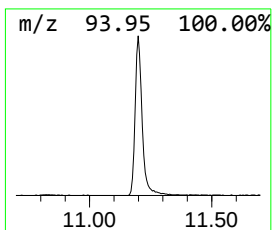
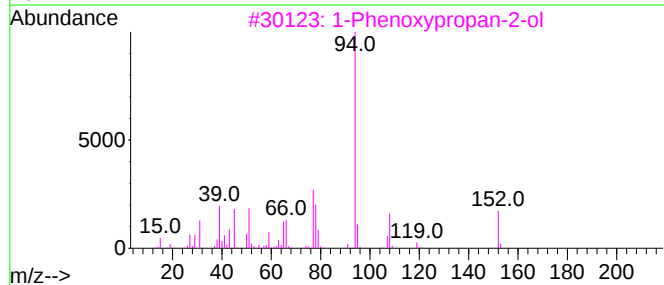
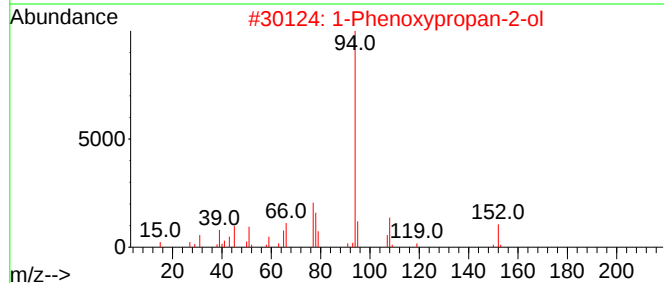
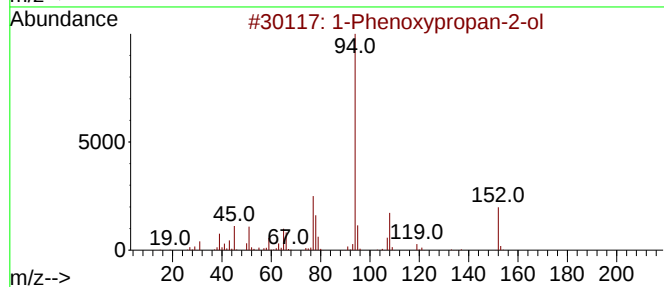
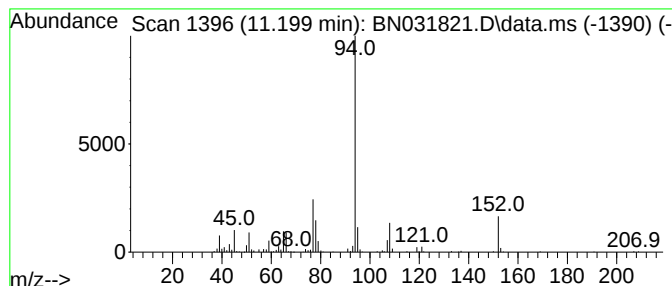
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Quant Title : SVOA CALIBRATION

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

Peak Number 3 1-Phenoxypropan-2-ol Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.199	2.28 ng/ul	359359	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	96
3			1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	94
4			1-Propanol, 3-phenoxy-	152	C9H12O2	006180-61-6	87
5			1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN060624\
Data File : BN031821.D
Acq On : 06 Jun 2024 16:34
Operator : MA/JU
Sample : P2634-07
Misc :
ALS Vial : 10 Sample Multiplier: 1

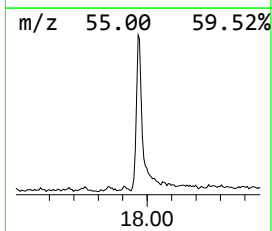
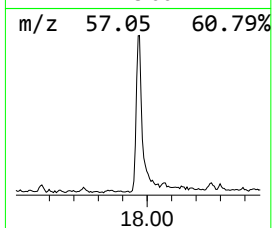
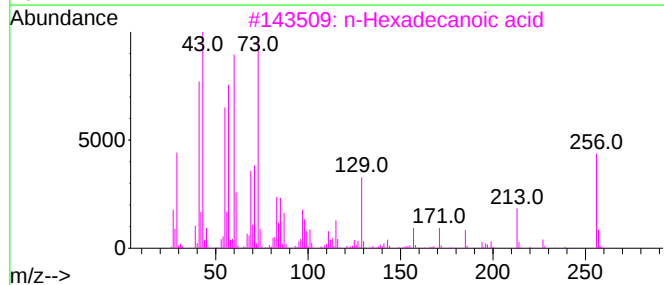
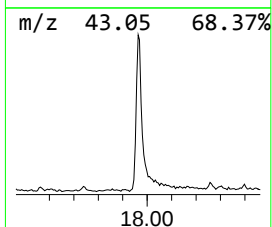
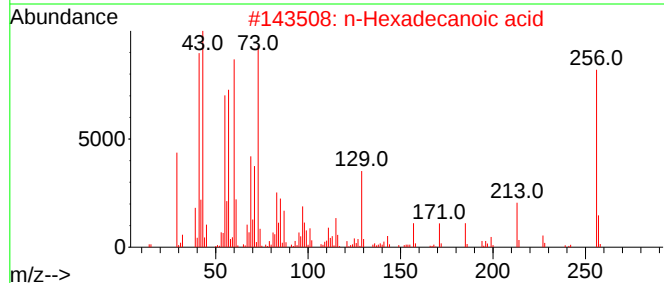
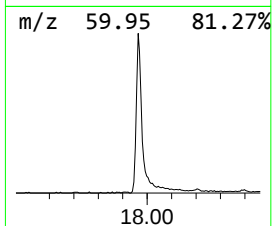
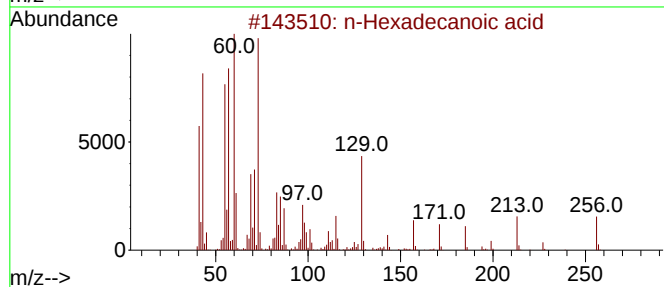
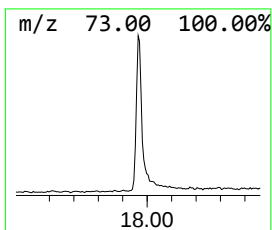
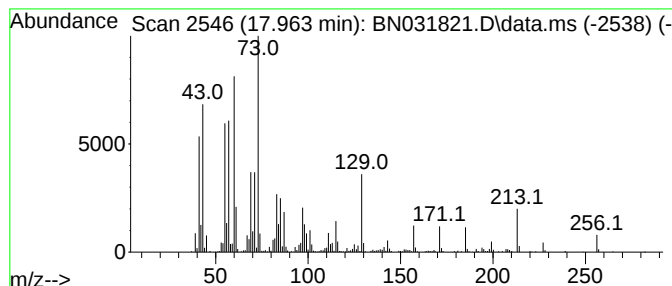
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ClientSampleId :
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 n-Hexadecanoic acid Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.	
17.963	4.08 ng/ul	983916	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	98
4	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
5	Tridecanoic acid	214	C13H26O2	000638-53-9	92



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN060624\
Data File : BN031821.D
Acq On : 06 Jun 2024 16:34
Operator : MA/JU
Sample : P2634-07
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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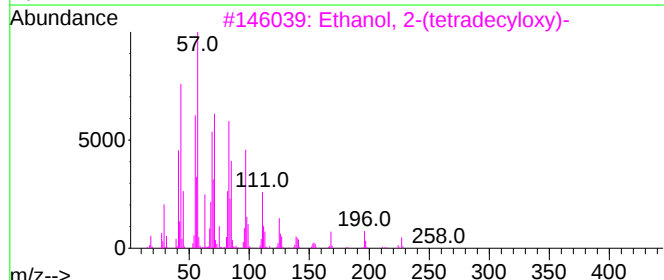
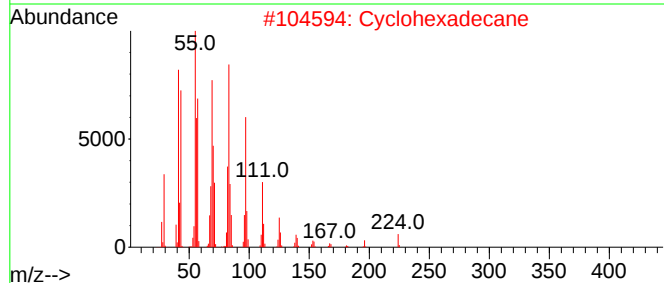
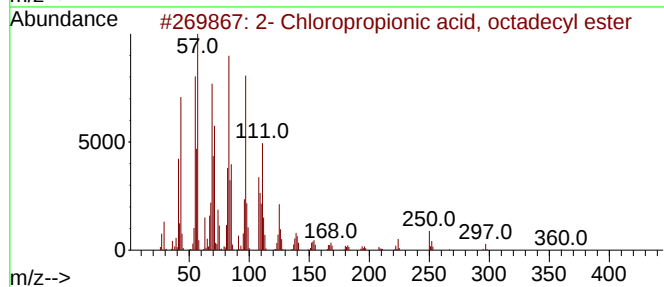
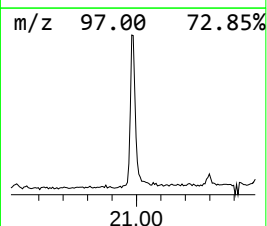
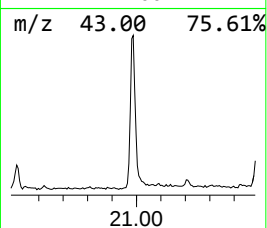
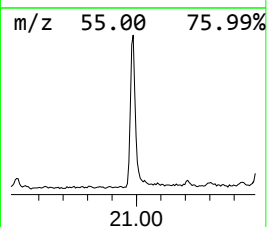
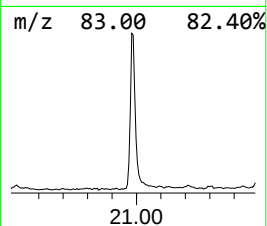
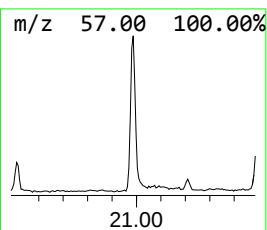
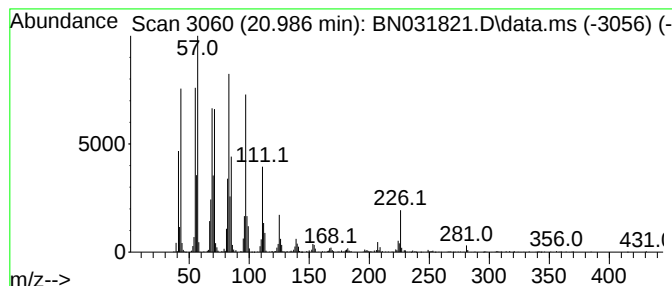
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 2- Chloropropionic acid, oc... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.986	3.96 ng/ul	1019280	Chrysene-d12	21.298

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2- Chloropropionic acid, octadec...	360	C21H41ClO2	088104-31-8	95
2			Cyclohexadecane	224	C16H32	000295-65-8	94
3			Ethanol, 2-(tetradecyloxy)-	258	C16H34O2	002136-70-1	94
4			1-Heneicosyl formate	340	C22H44O2	077899-03-7	93
5			Pentadecafluorooctanoic acid, oc...	666	C26H37F15O2	1000406-04-8	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN060624\
Data File : BN031821.D
Acq On : 06 Jun 2024 16:34
Operator : MA/JU
Sample : P2634-07
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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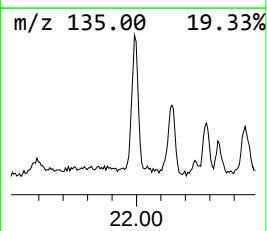
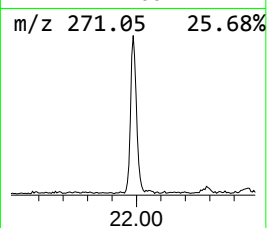
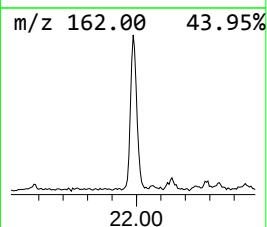
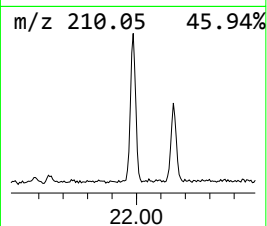
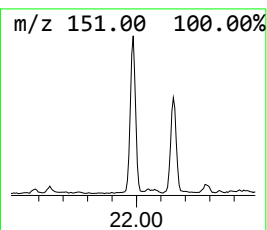
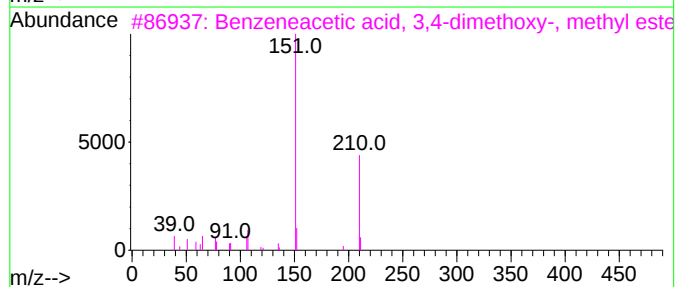
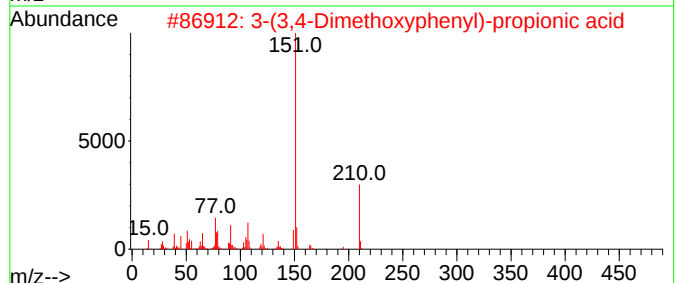
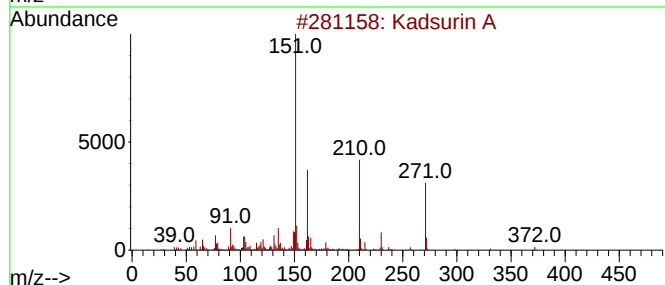
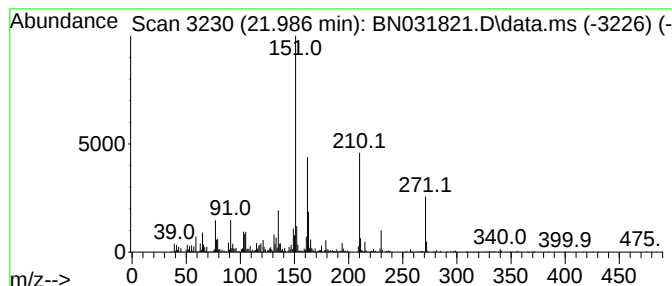
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 Kadsurin A Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.986	3.41 ng/ul	878897	Chrysene-d12	21.298

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Kadsurin A	372	C21H24O6	099340-07-5	87
2			3-(3,4-Dimethoxyphenyl)-propioni...	210	C11H14O4	002107-70-2	49
3			Benzeneacetic acid, 3,4-dimethox...	210	C11H14O4	015964-79-1	46
4			3-(3,4-Dimethoxyphenyl)-propioni...	210	C11H14O4	002107-70-2	46
5			Benzeneacetic acid, 3,4-dimethox...	210	C11H14O4	015964-79-1	43



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN060624\
Data File : BN031821.D
Acq On : 06 Jun 2024 16:34
Operator : MA/JU
Sample : P2634-07
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
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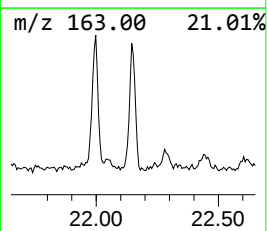
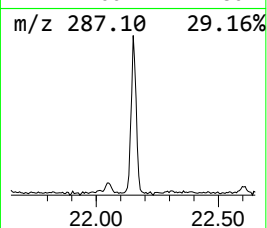
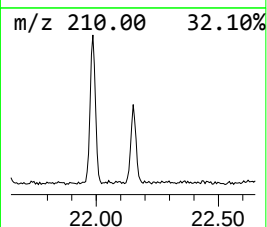
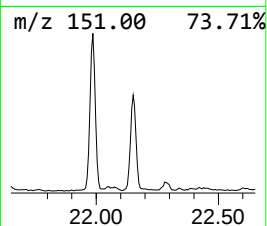
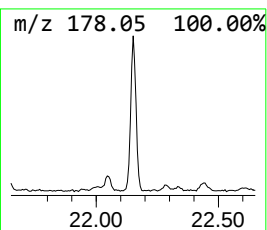
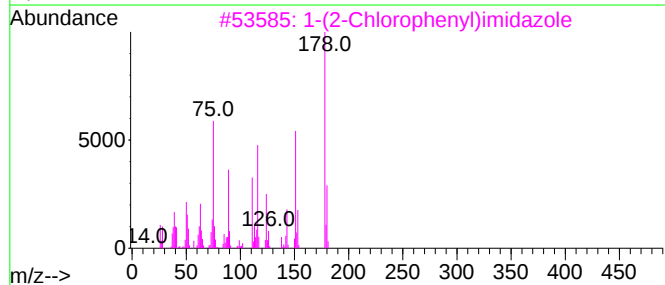
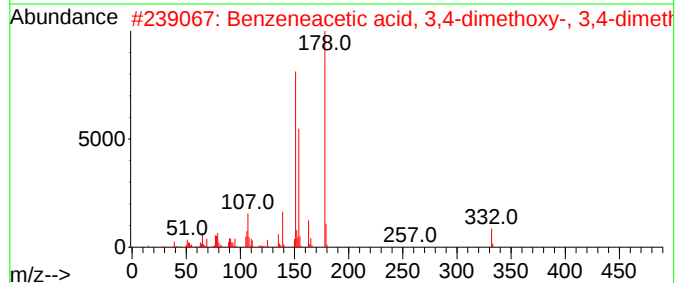
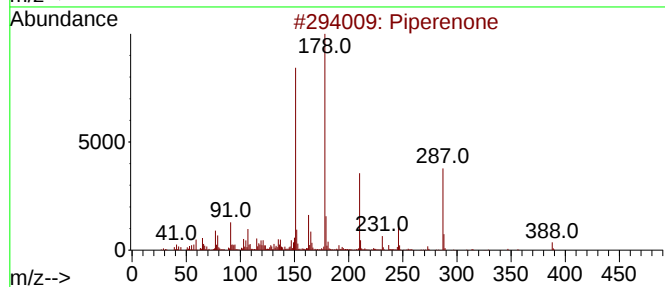
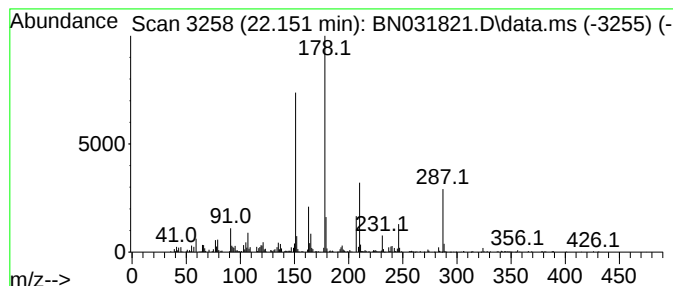
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 Piperone Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.151	2.57 ng/ul	662617	Chrysene-d12	21.298

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Piperone	388	C22H28O6	057625-31-7	87
2			Benzeneacetic acid, 3,4-dimethox...	332	C18H20O6	024195-27-5	53
3			1-(2-Chlorophenyl)imidazole	178	C9H7ClN2	051581-50-1	47
4			Methyleugenol	178	C11H14O2	000093-15-2	46
5			7-Isoquinolinol, 1,2,3,4-tetrahy...	285	C17H19NO3	024656-61-9	42



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN060624\
Data File : BN031821.D
Acq On : 06 Jun 2024 16:34
Operator : MA/JU
Sample : P2634-07
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
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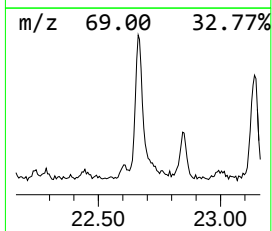
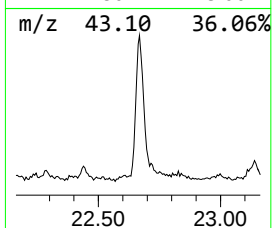
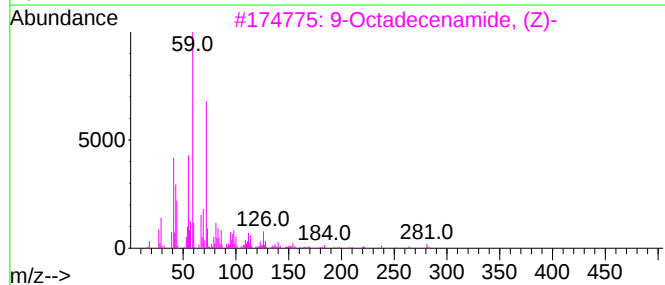
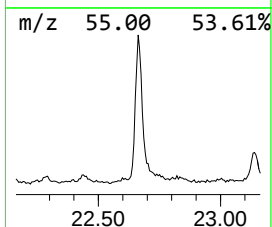
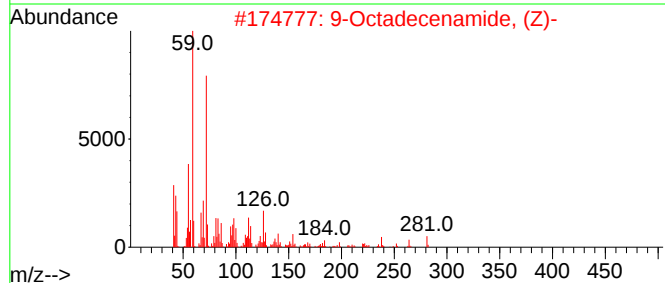
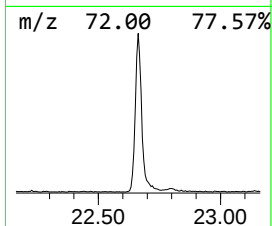
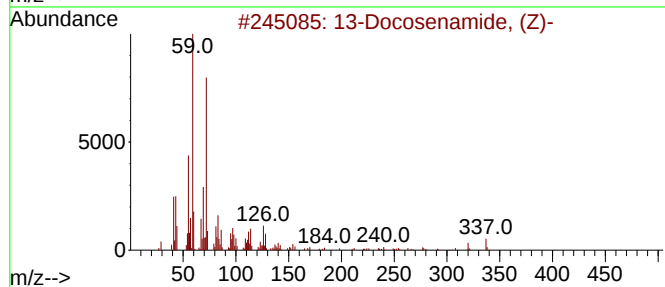
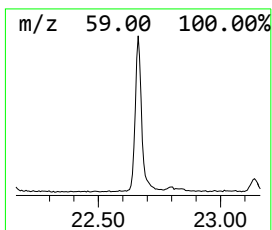
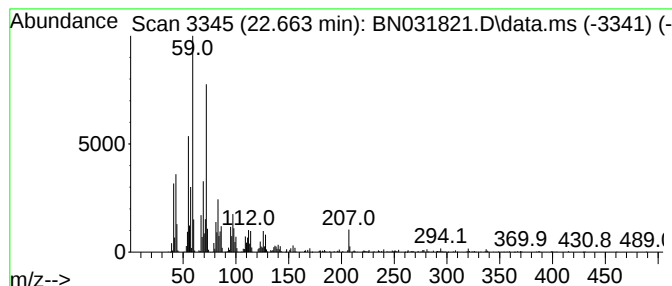
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 13-Docosenamide, (Z)- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.663	3.35 ng/ul	862449	Chrysene-d12	21.298

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			13-Docosenamide, (Z)-	337	C22H43NO	000112-84-5	99
2			9-Octadecenamide, (Z)-	281	C18H35NO	000301-02-0	93
3			9-Octadecenamide, (Z)-	281	C18H35NO	000301-02-0	86
4			Palmitoleamide	253	C16H31NO	106010-22-4	80
5			13-Docosenamide, (Z)-	337	C22H43NO	000112-84-5	50



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN060624\
Data File : BN031821.D
Acq On : 06 Jun 2024 16:34
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Sample : P2634-07
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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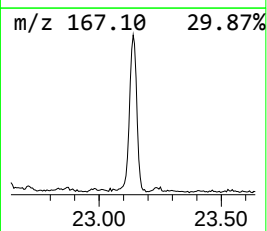
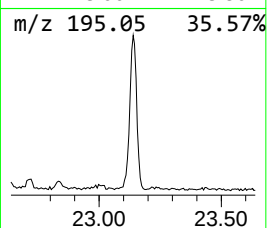
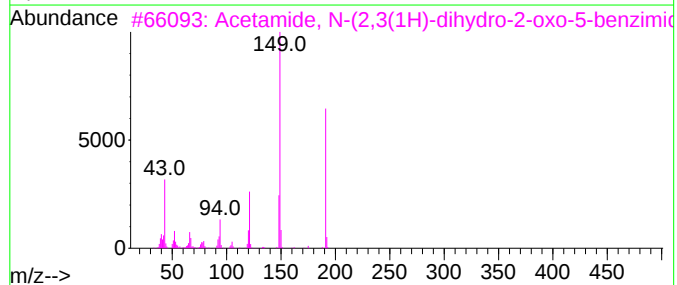
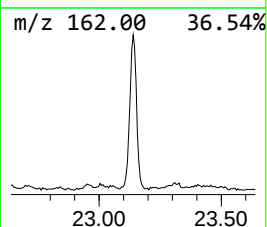
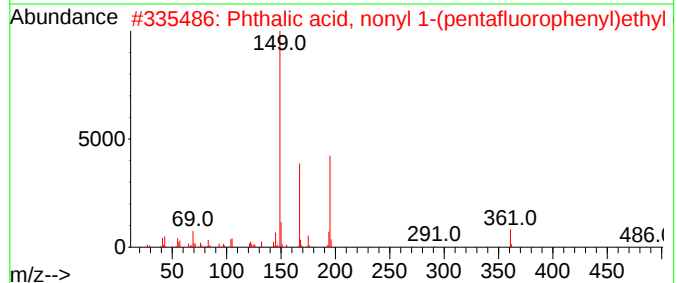
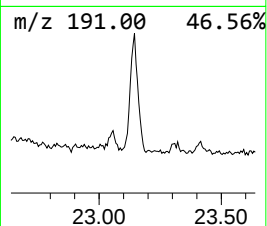
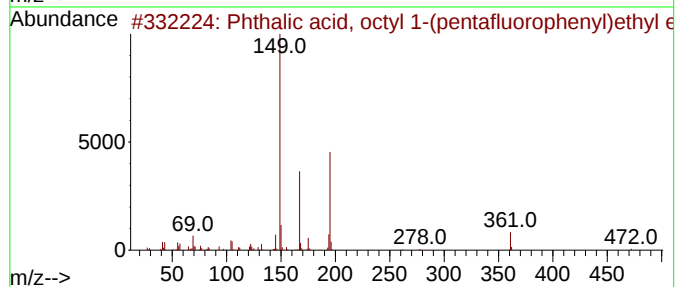
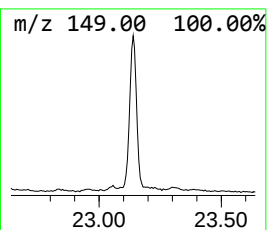
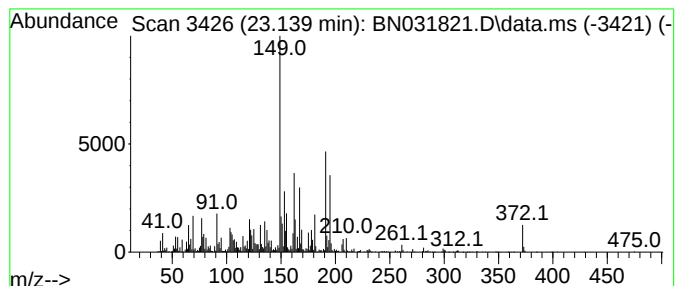
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 unknown-01 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.139	10.96 ng/ul	2166520	Perylene-d12	24.233

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phthalic acid, octyl 1-(pentaflu...	472	C24H25F5O4	1000382-93-1	43
2			Phthalic acid, nonyl 1-(pentaflu...	486	C25H27F5O4	1000382-93-2	43
3			Acetamide, N-(2,3(1H)-dihydro-2-...	191	C9H9N3O2	091085-68-6	38
4			Phthalic acid, 3,5-dinitro-4-met...	458	C23H26N2O8	1000415-51-5	35
5			Mono(2-ethylhexyl) phthalate	278	C16H22O4	004376-20-9	35



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN060624\
Data File : BN031821.D
Acq On : 06 Jun 2024 16:34
Operator : MA/JU
Sample : P2634-07
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
BHBX1

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

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
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.beta.-Pinene	7.105	2.4	ng/ul	251960	1	7.669	2141140	20.0
1-Phenoxypropan...	11.199	2.3	ng/ul	359359	2	10.451	3148040	20.0
n-Hexadecanoic ...	17.963	4.1	ng/ul	983916	4	17.063	4827720	20.0
2- Chloropropio...	20.986	4.0	ng/ul	1019280	5	21.298	5151380	20.0
Kadsurin A	21.986	3.4	ng/ul	878897	5	21.298	5151380	20.0
Piperenone	22.151	2.6	ng/ul	662617	5	21.298	5151380	20.0
13-Docosenamide...	22.663	3.4	ng/ul	862449	5	21.298	5151380	20.0
unknown-01	23.139	11.0	ng/ul	2166520	6	24.233	3953950	20.0