

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN032923\
 Data File : BN024602.D
 Acq On : 29 Mar 2023 10:03
 Operator : CG/JU
 Sample : 01982-01
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 EYGB1

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-BN032923.M
 Title : SVOA CALIBRATION

Signal : TIC: BN024602.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.364	9	13	25	rVB	66341	108177	1.65%	0.127%
2	3.793	83	86	119	rBV	825186	1412648	21.61%	1.658%
3	4.028	123	126	131	rVB5	10061	14295	0.22%	0.017%
4	4.128	139	143	149	rBV	25873	35068	0.54%	0.041%
5	4.234	158	161	166	rVB6	6589	11682	0.18%	0.014%
6	4.546	213	214	220	rVB6	4828	6917	0.11%	0.008%
7	5.064	296	302	310	rVB	23292	41406	0.63%	0.049%
8	7.152	649	657	670	rVB	1268282	1991868	30.47%	2.338%
9	7.322	679	686	698	rVB	953342	1468120	22.46%	1.723%
10	7.522	712	720	730	rBV	1163411	1825115	27.92%	2.142%
11	7.999	794	801	807	rBV	1293455	2054663	31.43%	2.411%
12	8.058	809	811	817	rVB7	8184	12203	0.19%	0.014%
13	8.222	836	839	843	rBV5	9178	13813	0.21%	0.016%
14	8.534	888	892	895	rBV5	9166	13242	0.20%	0.016%
15	8.705	910	921	932	rBV	1524703	2362030	36.13%	2.772%
16	8.875	945	950	953	rBV5	5750	8611	0.13%	0.010%
17	9.163	992	999	1009	rBV	1061395	1765388	27.01%	2.072%
18	9.328	1022	1027	1031	rBV5	10598	19085	0.29%	0.022%
19	9.522	1055	1060	1064	rBV8	3334	7213	0.11%	0.008%
20	9.575	1064	1069	1075	rVB3	20777	35998	0.55%	0.042%
21	9.893	1115	1123	1133	rBV	890813	1428072	21.85%	1.676%
22	10.210	1168	1177	1182	rVB3	11716	24491	0.37%	0.029%
23	10.281	1185	1189	1193	rBV3	9400	13611	0.21%	0.016%
24	10.428	1206	1214	1228	rBV	1390345	2263916	34.63%	2.657%
25	10.816	1268	1280	1288	rBV	1836700	3096259	47.37%	3.634%
26	10.952	1290	1303	1318	rVV	1363753	2277783	34.84%	2.673%
27	11.599	1406	1413	1417	rBV7	8120	15282	0.23%	0.018%
28	11.834	1449	1453	1461	rVB9	4565	9501	0.15%	0.011%
29	12.222	1513	1519	1523	rBV3	13429	20001	0.31%	0.023%
30	12.310	1527	1534	1537	rBV9	3729	7680	0.12%	0.009%
31	12.399	1543	1549	1554	rVB	23379	36591	0.56%	0.043%
32	12.846	1619	1625	1631	rBV9	5295	11582	0.18%	0.014%
33	12.981	1642	1648	1654	rBV6	12218	31571	0.48%	0.037%
34	13.246	1689	1693	1697	rVB6	4608	7085	0.11%	0.008%
35	13.457	1724	1729	1734	rVB3	19217	28541	0.44%	0.033%
36	13.610	1750	1755	1760	rVB7	5647	10892	0.17%	0.013%

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Integrator: RTE
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37	13.898	1799	1804	1810	rBV6	9522	19749	0.30%	0.023%
38	14.045	1822	1829	1836	rBV	2337224	3153499	48.24%	3.701%
39	14.157	1845	1848	1853	rVB5	6567	10029	0.15%	0.012%
40	14.334	1871	1878	1889	rVV	2547219	3706601	56.70%	4.350%
41	14.522	1906	1910	1915	rVB6	11912	17335	0.27%	0.020%
42	14.640	1923	1930	1939	rBV	2803108	4017938	61.47%	4.715%
43	14.828	1955	1962	1979	rBV	1145555	1612797	24.67%	1.893%
44	15.051	1994	2000	2006	rBV5	23409	41816	0.64%	0.049%
45	15.410	2057	2061	2062	rBV3	9857	8454	0.13%	0.010%
46	15.628	2090	2098	2106	rBV	3357368	4828545	73.87%	5.667%
47	15.745	2112	2118	2128	rBV	965452	1304737	19.96%	1.531%
48	15.869	2134	2139	2145	rBV3	14593	23749	0.36%	0.028%
49	15.992	2155	2160	2164	rBV	60741	84407	1.29%	0.099%
50	16.334	2216	2218	2225	rVB8	7670	11519	0.18%	0.014%
51	16.428	2229	2234	2237	rVV7	5613	10322	0.16%	0.012%
52	16.498	2242	2246	2255	rVV9	13790	27549	0.42%	0.032%
53	16.763	2289	2291	2293	rBV3	7709	8462	0.13%	0.010%
54	16.845	2303	2305	2309	rBV5	6571	8378	0.13%	0.010%
55	16.875	2309	2310	2315	rVB4	8168	6563	0.10%	0.008%
56	17.069	2338	2343	2348	rVB7	5245	9952	0.15%	0.012%
57	17.134	2348	2354	2357	rBV8	11672	20493	0.31%	0.024%
58	17.169	2357	2360	2364	rVB4	9764	16310	0.25%	0.019%
59	17.292	2379	2381	2386	rVB6	6117	7204	0.11%	0.008%
60	17.386	2390	2397	2408	rBV	3450288	4744630	72.58%	5.568%
61	17.486	2408	2414	2425	rVV	3617211	5148619	78.76%	6.042%
62	17.575	2425	2429	2436	rVB5	39211	59824	0.92%	0.070%
63	17.745	2454	2458	2459	rBV4	12934	17738	0.27%	0.021%
64	18.045	2506	2509	2513	rVB6	7734	10354	0.16%	0.012%
65	18.269	2543	2547	2556	rBV	252348	341059	5.22%	0.400%
66	18.345	2557	2560	2566	rVB	36996	50913	0.78%	0.060%
67	18.698	2616	2620	2624	rVB2	60505	73204	1.12%	0.086%
68	19.339	2725	2729	2736	rBV2	43510	67808	1.04%	0.080%
69	19.410	2737	2741	2750	rVV2	50704	105931	1.62%	0.124%
70	19.539	2760	2763	2769	rBV3	61805	80318	1.23%	0.094%
71	19.698	2787	2790	2794	rBV	38571	46414	0.71%	0.054%
72	19.775	2797	2803	2815	rVV	4628524	6061862	92.73%	7.114%
73	20.310	2892	2894	2897	rVB2	31369	24893	0.38%	0.029%
74	21.239	3046	3052	3055	rBV	1730439	2777439	42.49%	3.259%
75	21.292	3055	3061	3067	rVV	2640610	4142354	63.37%	4.861%
76	21.345	3067	3070	3078	rVV	814161	1025333	15.69%	1.203%

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77	21.416	3078	3082	3086	rVV	431503	523554	8.01%	0.614%
78	21.480	3090	3093	3100	rVB	461993	533694	8.16%	0.626%
79	21.574	3104	3109	3116	rVV2	3822556	5229330	80.00%	6.137%
80	23.286	3397	3400	3408	rVB2	192237	304874	4.66%	0.358%
81	23.910	3499	3506	3522	rVB2	3164489	6536918	100.00%	7.671%
82	24.068	3526	3533	3542	rVB2	2391285	5314524	81.30%	6.237%
83	24.692	3635	3639	3646	rVB2	278226	542658	8.30%	0.637%

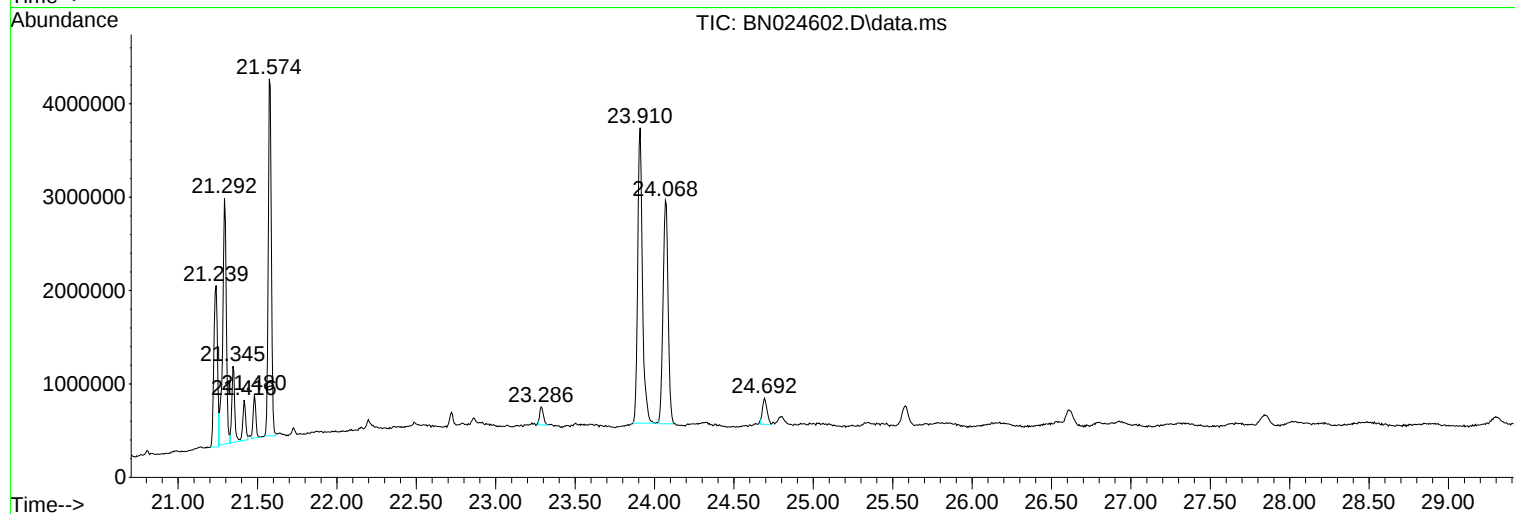
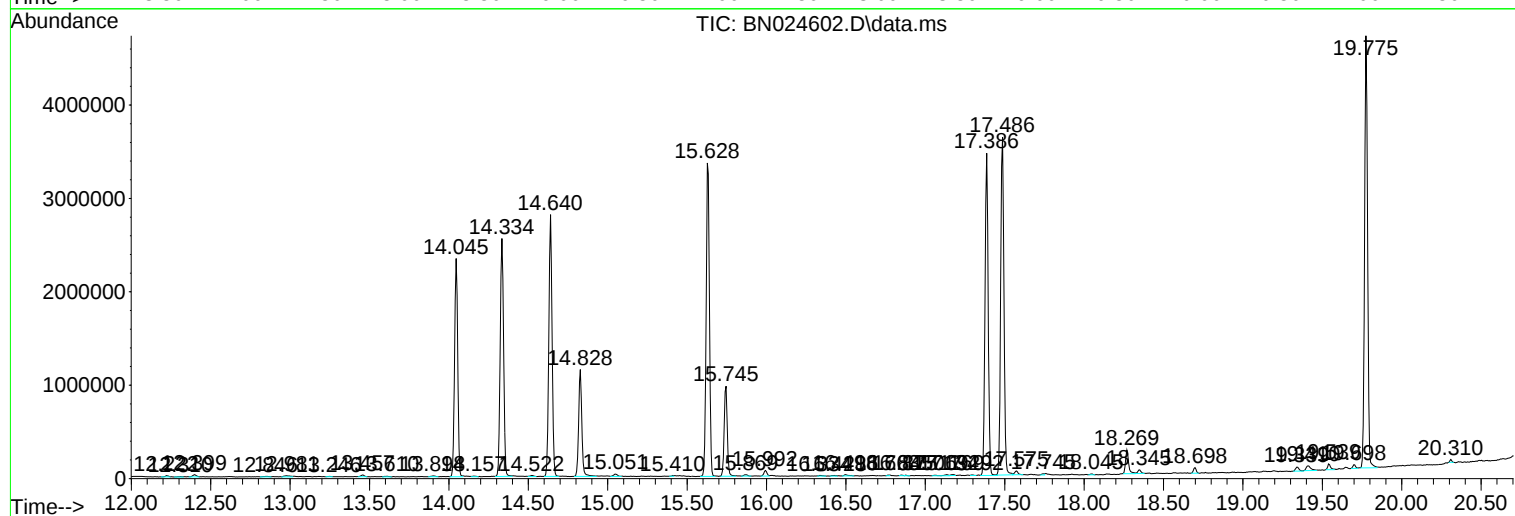
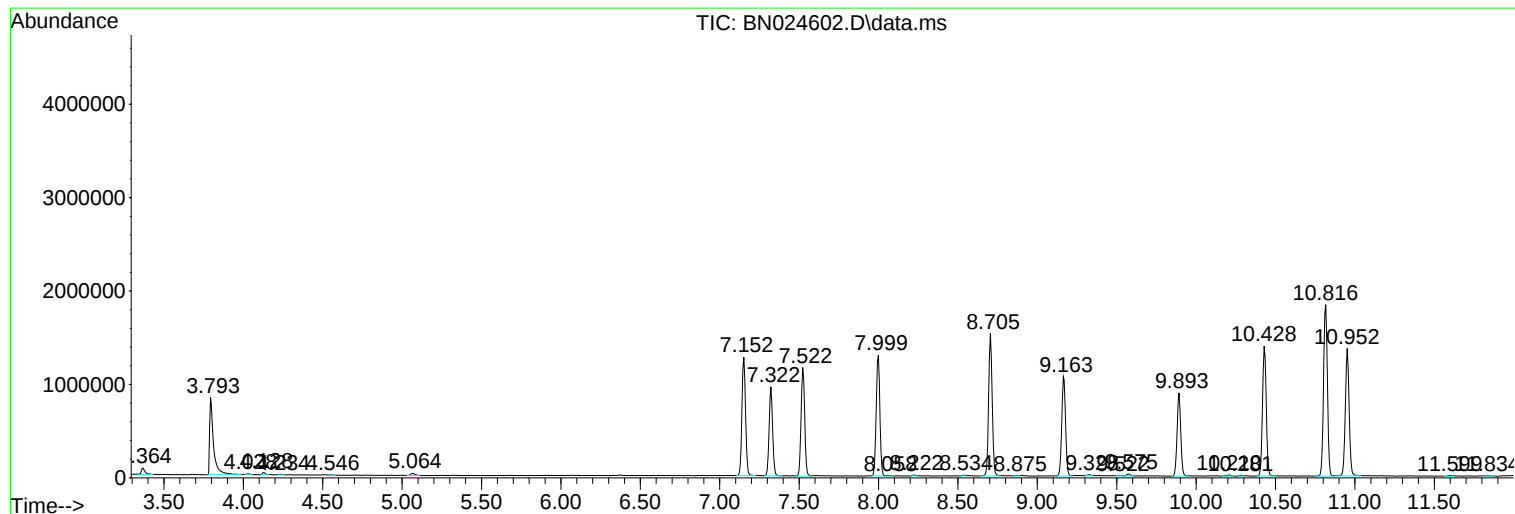
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Instrument :
BNA_N
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN032923.M
Quant Title : SVOA CALIBRATION

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



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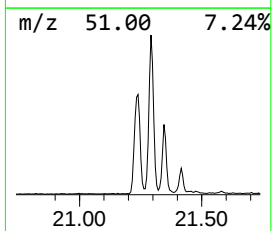
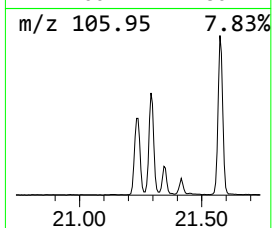
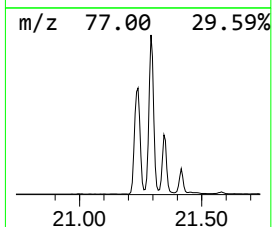
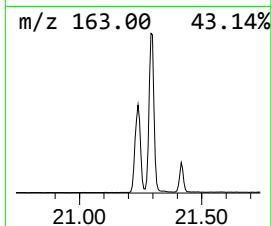
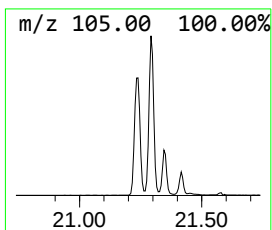
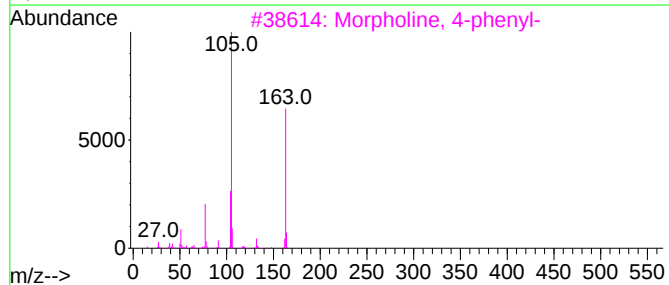
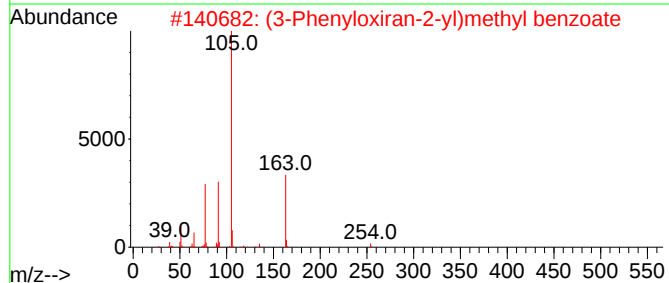
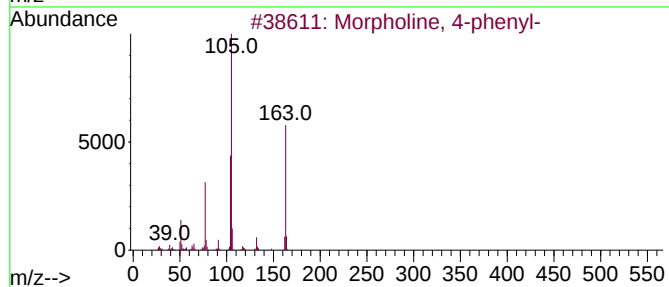
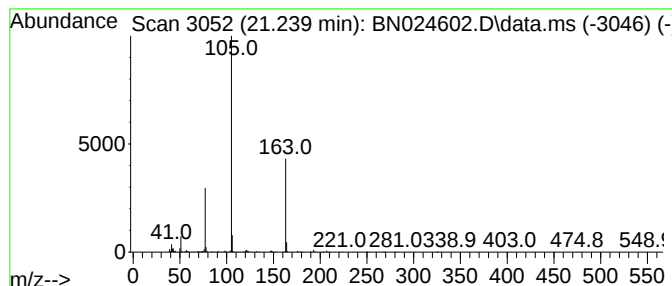
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN032923.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Morpholine, 4-phenyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.239	10.62 ng/ul	2777440	Chrysene-d12	21.580

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Morpholine, 4-phenyl-	163	C10H13NO	000092-53-5	72
2			(3-Phenyloxiran-2-yl)methyl benz...	254	C16H14O3	1000366-96-1	53
3			Morpholine, 4-phenyl-	163	C10H13NO	000092-53-5	53
4			Oxybis(propane-1,2-diyl) dibenzoate	342	C20H22O5	000094-03-1	43
5			Isopropyl phenyl ketone	148	C10H12O	000611-70-1	38



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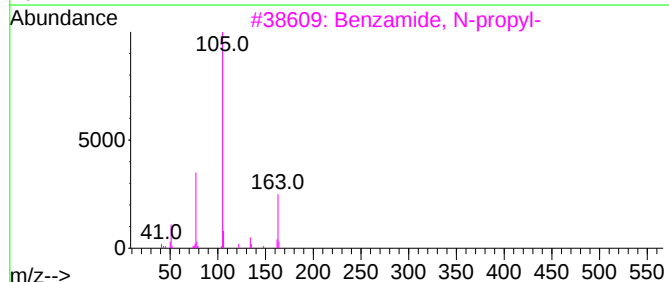
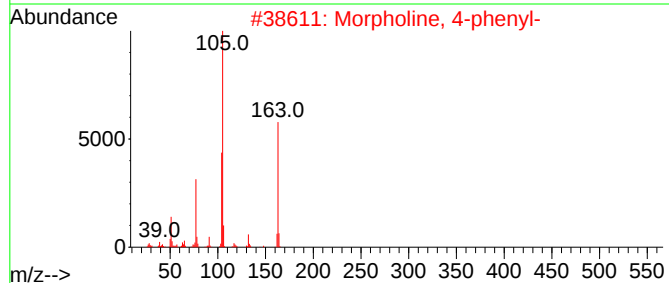
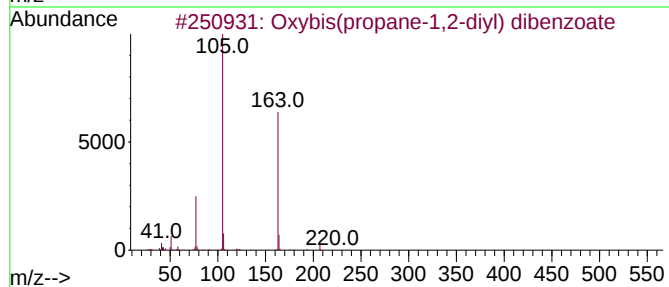
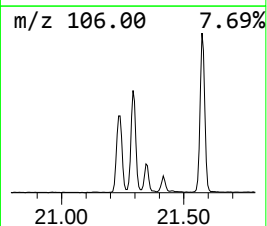
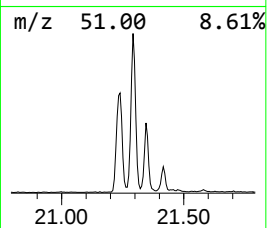
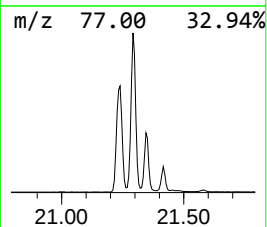
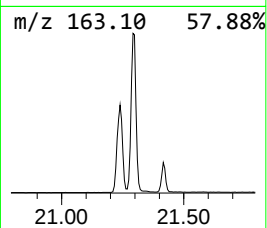
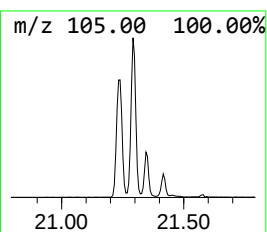
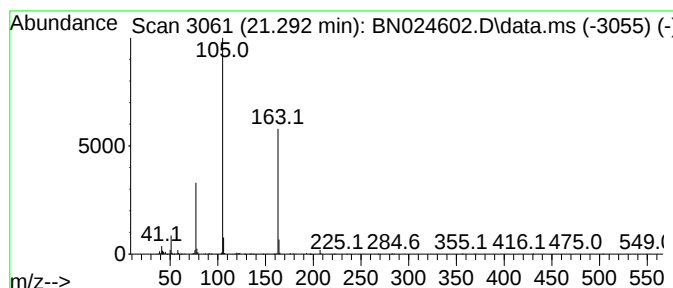
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 Oxybis(propene-1,2-diyl) di... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.292	15.84 ng/ul	4142350	Chrysene-d12	21.580

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Oxybis(propene-1,2-diyl) dibenzoate	342	C20H22O5	000094-03-1	90
2			Morpholine, 4-phenyl-	163	C10H13NO	000092-53-5	72
3			Benzamide, N-propyl-	163	C10H13NO	010546-70-0	72
4			Morpholine, 4-phenyl-	163	C10H13NO	000092-53-5	64
5			Morpholine, 4-phenyl-	163	C10H13NO	000092-53-5	64



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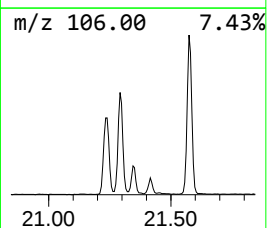
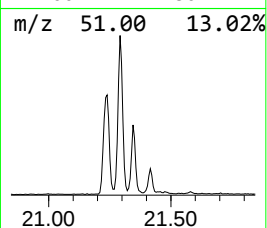
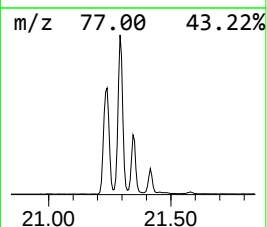
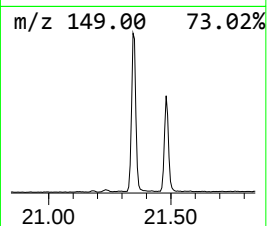
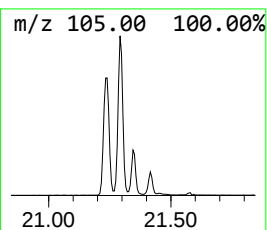
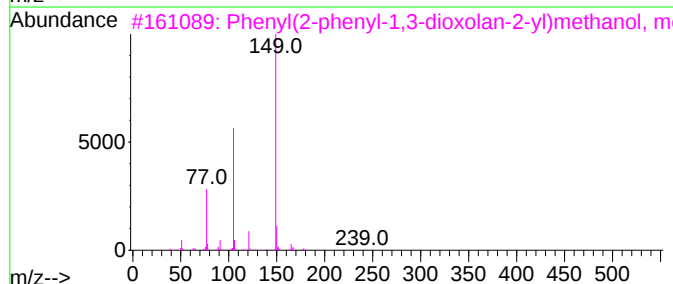
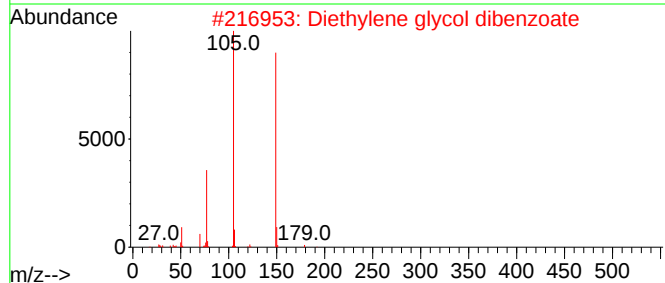
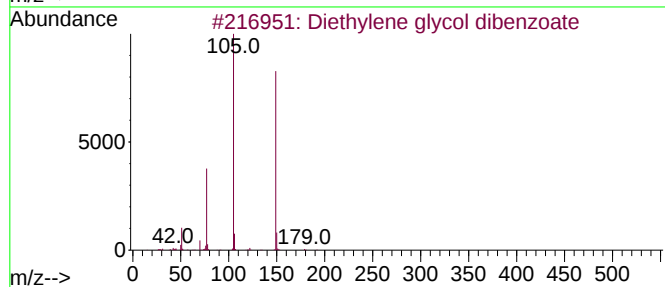
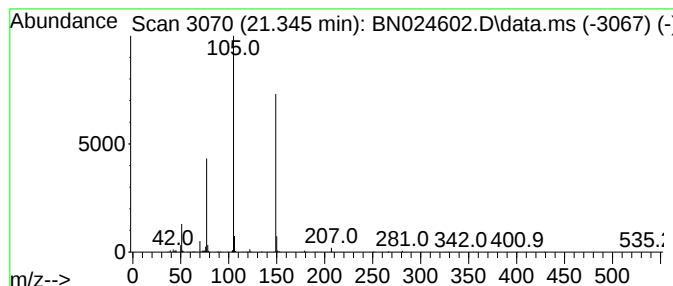
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Quant Title : SVOA CALIBRATION

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

Peak Number 3 Diethylene glycol dibenzoate Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.345	3.92 ng/ul	1025330	Chrysene-d12	21.580

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Diethylene glycol dibenzoate	314	C18H18O5	000120-55-8	90
2			Diethylene glycol dibenzoate	314	C18H18O5	000120-55-8	72
3			Phenyl(2-phenyl-1,3-dioxolan-2-y...	270	C17H18O3	1000507-07-6	64
4			Ethanone, 2-hydroxy-1-phenyl-	136	C8H8O2	000582-24-1	47
5			N-(5-Pyrimidinyl)benzamide, Ac d...	241	C13H11N3O2	1000504-27-8	47



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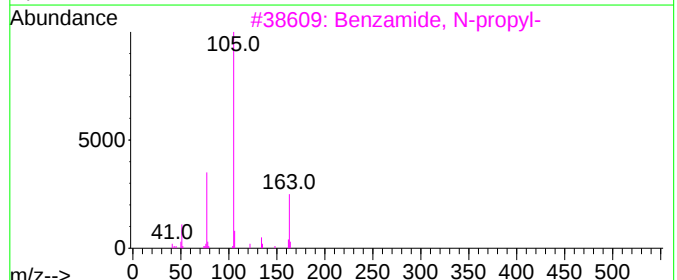
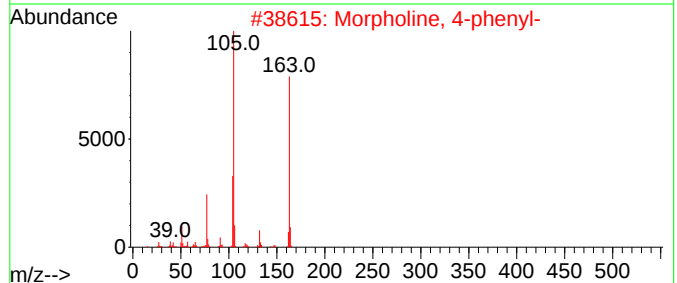
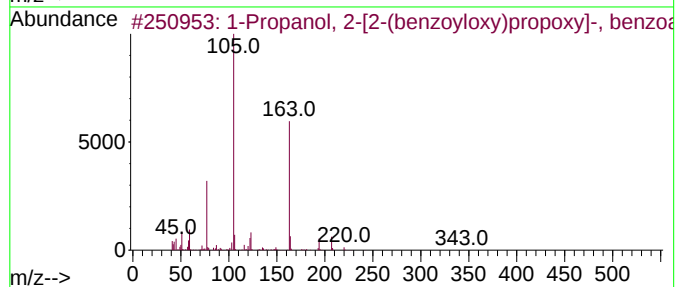
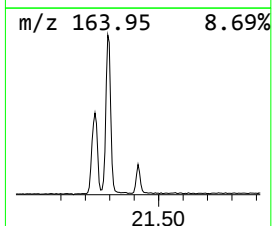
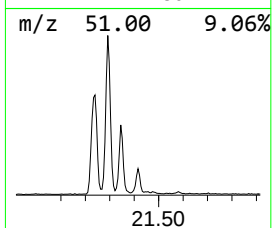
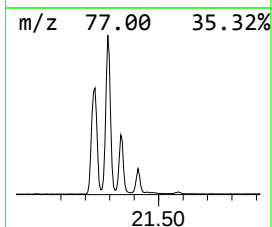
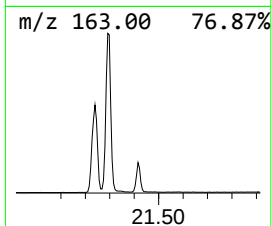
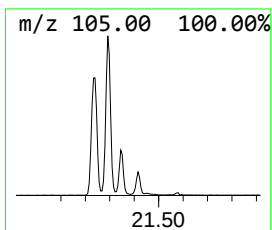
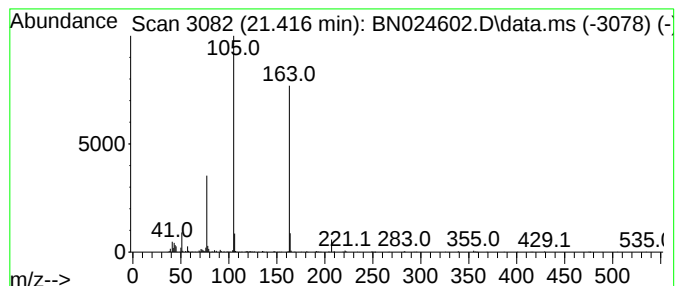
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 1-Propanol, 2-[2-(benzoylox... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.416	2.00 ng/ul	523554	Chrysene-d12	21.580

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Propanol, 2-[2-(benzoyloxy)pro...	342	C20H22O5	020109-39-1	74
2			Morpholine, 4-phenyl-	163	C10H13NO	000092-53-5	72
3			Benzamide, N-propyl-	163	C10H13NO	010546-70-0	72
4			Morpholine, 4-phenyl-	163	C10H13NO	000092-53-5	72
5			Glyoxylamide, N-phenyl-	163	C9H9NO2	032331-52-5	59



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN032923\
Data File : BN024602.D
Acq On : 29 Mar 2023 10:03
Operator : CG/JU
Sample : 01982-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

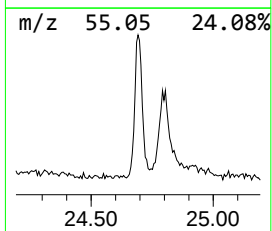
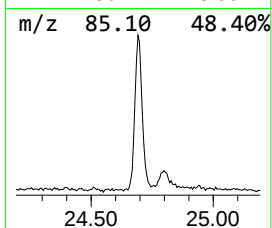
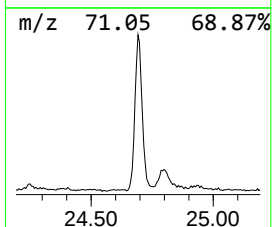
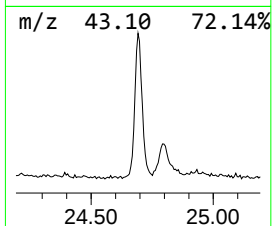
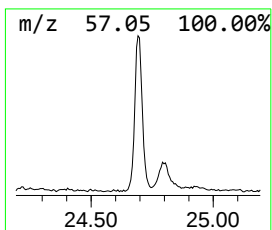
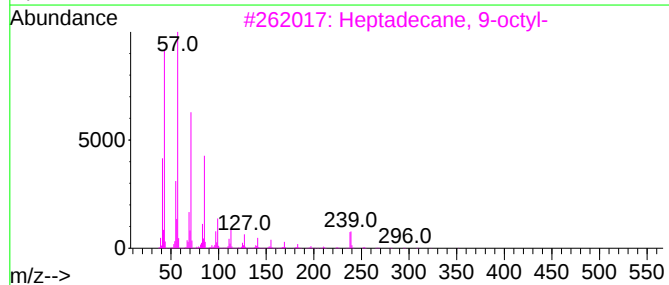
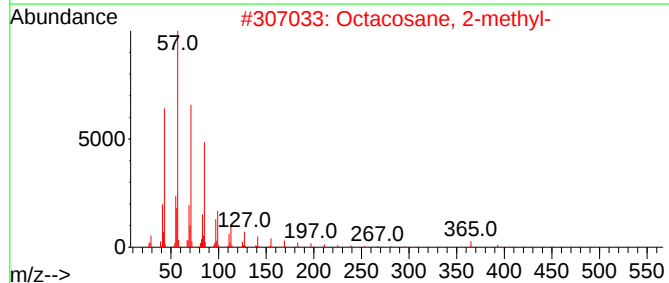
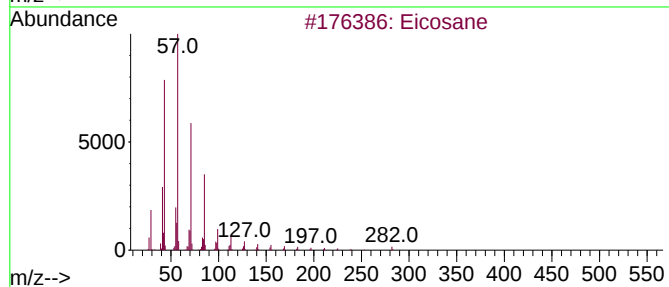
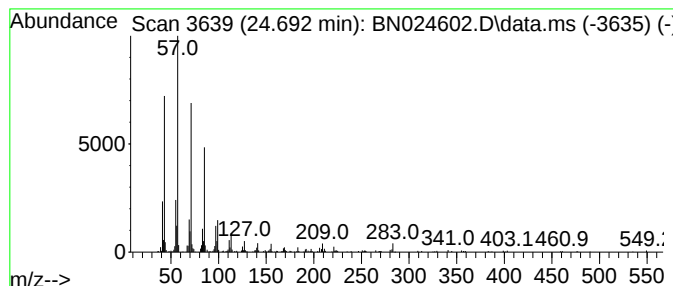
Instrument :
BNA_N
ClientSampleId :
EYGB1

Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN032923.M
Quant Title : SVOA CALIBRATION

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

Peak Number 5 (DEL) Alkane: Straight-Chai... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD		R.T.	
24.692	2.04 ng/ul	542658	Perylene-d12		24.069	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Eicosane		282	C20H42	000112-95-8	95
2	Octacosane, 2-methyl-		408	C29H60	001560-98-1	90
3	Heptadecane, 9-octyl-		352	C25H52	007225-64-1	90
4	Hexacosane		366	C26H54	000630-01-3	90
5	Octacosane		394	C28H58	000630-02-4	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN032923\
Data File : BN024602.D
Acq On : 29 Mar 2023 10:03
Operator : CG/JU
Sample : 01982-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
EYGB1

Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN032923.M
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
Morpholine, 4-p...	21.239	10.6	ng/ul	2777440	5	21.580	5229330	20.0
Oxybis(propane-...	21.292	15.8	ng/ul	4142350	5	21.580	5229330	20.0
Diethylene glyc...	21.345	3.9	ng/ul	1025330	5	21.580	5229330	20.0
1-Propanol, 2-[...	21.416	2.0	ng/ul	523554	5	21.580	5229330	20.0
(DEL) Alkane: S...	24.692	2.0	ng/ul	542658	6	24.069	5314520	20.0