

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN062424\
 Data File : BN032212.D
 Acq On : 24 Jun 2024 17:41
 Operator : MA/JU
 Sample : P2856-09
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 C0SH1

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: BN032212.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.169	26	31	36	rBV	49336	70461	1.36%	0.124%
2	3.434	73	76	88	rVB	30957	52312	1.01%	0.092%
3	3.569	96	99	120	rBV	630418	986958	19.05%	1.742%
4	3.881	147	152	157	rVB	15655	24813	0.48%	0.044%
5	4.375	233	236	241	rBV7	5625	10548	0.20%	0.019%
6	4.422	241	244	249	rVB4	6489	9496	0.18%	0.017%
7	4.769	298	303	310	rBV	24532	38804	0.75%	0.068%
8	5.163	365	370	375	rVB4	6555	9600	0.19%	0.017%
9	5.299	388	393	395	rBV4	6126	8306	0.16%	0.015%
10	5.658	449	454	459	rBV6	6423	10371	0.20%	0.018%
11	5.734	462	467	475	rBV	31945	55018	1.06%	0.097%
12	6.263	552	557	564	rVB10	4584	8949	0.17%	0.016%
13	6.722	631	635	638	rVB6	4237	5576	0.11%	0.010%
14	6.816	642	651	668	rVV	781480	1353345	26.12%	2.389%
15	6.975	672	678	690	rVB	671652	1066682	20.59%	1.883%
16	7.163	703	710	720	rVV	874525	1436521	27.73%	2.536%
17	7.246	720	724	726	rVV3	19798	32467	0.63%	0.057%
18	7.269	726	728	735	rVV3	22652	40278	0.78%	0.071%
19	7.628	776	789	797	rBV	881185	1426902	27.54%	2.519%
20	7.704	797	802	814	rVB4	25290	70207	1.36%	0.124%
21	7.993	845	851	855	rBV5	3226	5534	0.11%	0.010%
22	8.157	874	879	886	rBV	33212	54979	1.06%	0.097%
23	8.340	903	910	925	rBV	1028423	1740212	33.59%	3.072%
24	8.787	978	986	997	rBV	763205	1285760	24.82%	2.270%
25	8.940	1009	1012	1019	rVB7	5543	10231	0.20%	0.018%
26	9.163	1043	1050	1057	rBV4	7273	12948	0.25%	0.023%
27	9.351	1075	1082	1089	rBV	93264	150730	2.91%	0.266%
28	9.504	1101	1108	1121	rBV	609531	1077133	20.79%	1.901%
29	9.828	1157	1163	1170	rVB9	4897	13236	0.26%	0.023%
30	9.904	1170	1176	1181	rBV9	6823	12463	0.24%	0.022%
31	10.040	1191	1199	1218	rBV	954248	1711046	33.03%	3.020%
32	10.410	1253	1262	1267	rBV	1089540	1918803	37.04%	3.387%
33	10.457	1267	1270	1278	rVV	145043	227775	4.40%	0.402%
34	10.557	1278	1287	1305	rVV	955476	1753577	33.85%	3.095%
35	10.963	1350	1356	1363	rBV3	21447	43644	0.84%	0.077%
36	11.157	1383	1389	1408	rBV	87972	214809	4.15%	0.379%

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Integration Parameters: LSCINT.P

Integrator: RTE

Smoothing : OFF

Sampling : 1

Start Thrs: 0.2

Stop Thrs : 0

Filtering: 5

Min Area: 0.1 % of largest Peak

Max Peaks: 100

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	11.916	1514	1518	1523	rVB5	4636	5960	0.12%	0.011%
38	12.004	1523	1533	1540	rBV2	19270	37471	0.72%	0.066%
39	12.075	1540	1545	1553	rVV	31891	53179	1.03%	0.094%
40	12.298	1578	1583	1589	rVB2	25897	45683	0.88%	0.081%

41	12.869	1677	1680	1688	rBV7	3087	6340	0.12%	0.011%
42	13.092	1713	1718	1721	rBV7	5035	8510	0.16%	0.015%
43	13.234	1737	1742	1747	rBV7	4403	7386	0.14%	0.013%
44	13.292	1747	1752	1756	rBV8	4096	6853	0.13%	0.012%
45	13.439	1772	1777	1778	rBV5	5410	6889	0.13%	0.012%

46	13.457	1778	1780	1784	rVB5	6177	6758	0.13%	0.012%
47	13.592	1798	1803	1808	rVB5	8450	14469	0.28%	0.026%
48	13.692	1813	1820	1835	rVV	1707195	2434244	46.99%	4.297%
49	13.828	1839	1843	1847	rVB6	4374	6663	0.13%	0.012%
50	13.969	1856	1867	1880	rBV	1985341	3105668	59.95%	5.482%

51	14.163	1894	1900	1905	rBV6	7038	16259	0.31%	0.029%
52	14.275	1911	1919	1927	rBV	1662670	2454255	47.37%	4.332%
53	14.510	1949	1959	1984	rVV	531134	1126809	21.75%	1.989%
54	14.704	1985	1992	1997	rVB9	13684	29644	0.57%	0.052%
55	14.892	2022	2024	2029	rVB5	5606	8154	0.16%	0.014%

56	15.069	2048	2054	2058	rBV6	7855	12959	0.25%	0.023%
57	15.128	2059	2064	2065	rBV4	4846	5864	0.11%	0.010%
58	15.269	2081	2088	2096	rBV	2469697	3728680	71.97%	6.582%
59	15.410	2106	2112	2124	rBV	544331	801700	15.47%	1.415%
60	15.516	2126	2130	2136	rVB6	15694	27925	0.54%	0.049%

61	15.739	2165	2168	2174	rVB7	4591	6627	0.13%	0.012%
62	15.828	2180	2183	2185	rBV4	4545	5523	0.11%	0.010%
63	15.992	2207	2211	2216	rVB7	7177	12349	0.24%	0.022%
64	16.057	2218	2222	2225	rVB6	4226	5547	0.11%	0.010%
65	16.157	2234	2239	2246	rBV8	4565	9890	0.19%	0.017%

66	16.439	2282	2287	2290	rBV7	6278	10277	0.20%	0.018%
67	16.580	2307	2311	2313	rVB4	3959	5375	0.10%	0.009%
68	16.786	2342	2346	2347	rBV4	4547	5439	0.10%	0.010%
69	16.810	2348	2350	2354	rVB4	6535	6646	0.13%	0.012%
70	16.951	2369	2374	2378	rBV7	5677	8554	0.17%	0.015%

71	17.028	2380	2387	2392	rBV	1989064	2860073	55.21%	5.048%
72	17.128	2397	2404	2415	rVV	2748828	4046797	78.11%	7.143%
73	17.227	2418	2421	2430	rVB9	12560	25576	0.49%	0.045%
74	17.698	2498	2501	2505	rVB5	10165	13300	0.26%	0.023%
75	17.927	2535	2540	2561	rBV	466187	857153	16.55%	1.513%

76	18.251	2593	2595	2598	rBV4	6094	8130	0.16%	0.014%
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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

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Title : SVOA CALIBRATION

77	18.986	2716	2720	2726	rVB4	41958	59474	1.15%	0.105%
78	19.063	2729	2733	2736	rBV	51287	81791	1.58%	0.144%
79	19.216	2754	2759	2772	rBV	268118	519308	10.02%	0.917%
80	19.427	2789	2795	2802	rBV	3393325	4672230	90.18%	8.247%
81	20.469	2970	2972	2976	rBV	56920	53696	1.04%	0.095%
82	20.945	3050	3053	3062	rVB	243916	358954	6.93%	0.634%
83	21.263	3101	3107	3113	rBV2	2214365	3247002	62.67%	5.731%
84	23.980	3561	3569	3580	rVB2	2196720	5180721	100.00%	9.145%
85	24.174	3594	3602	3614	rVB	1485148	3684218	71.11%	6.503%

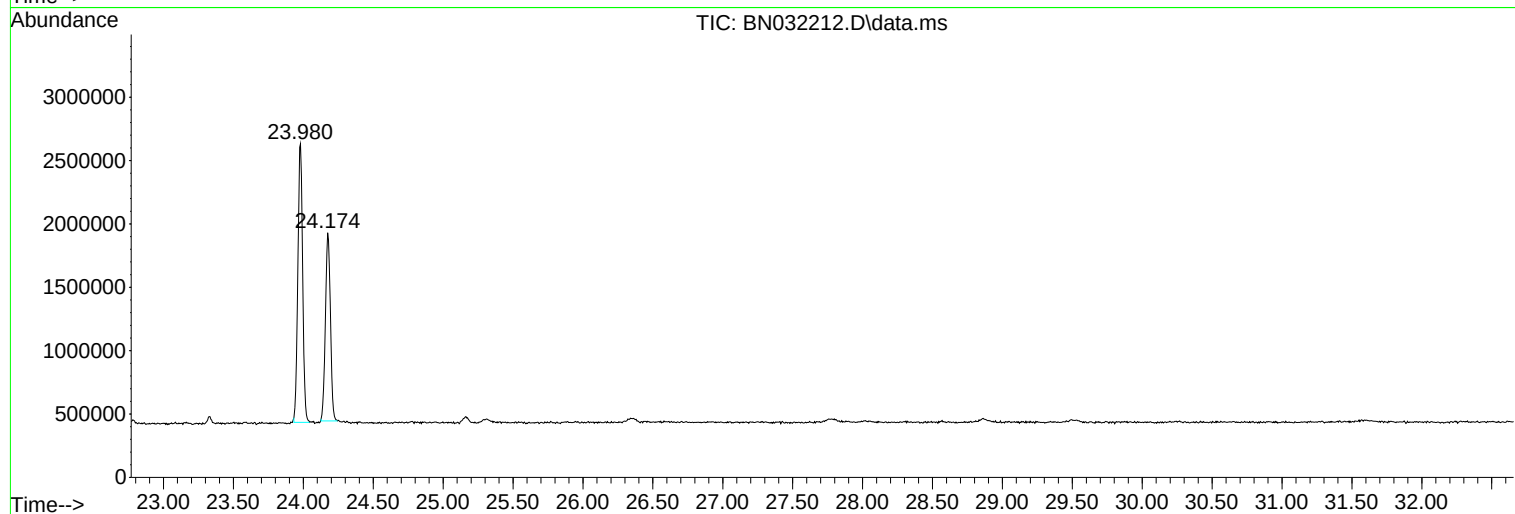
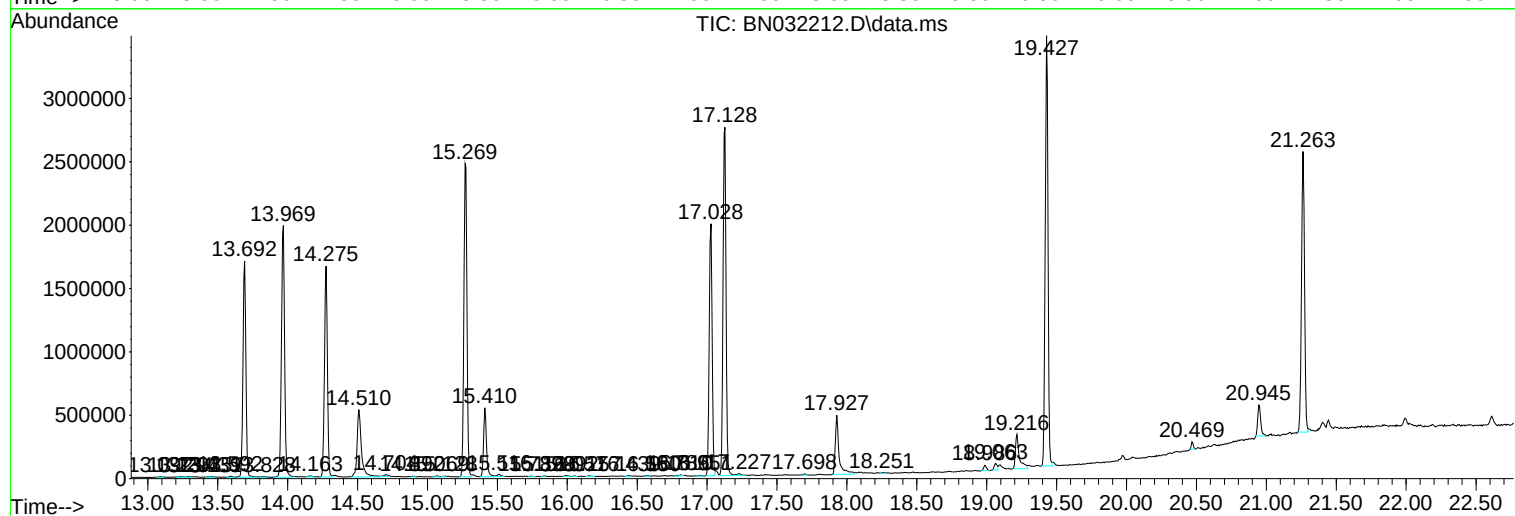
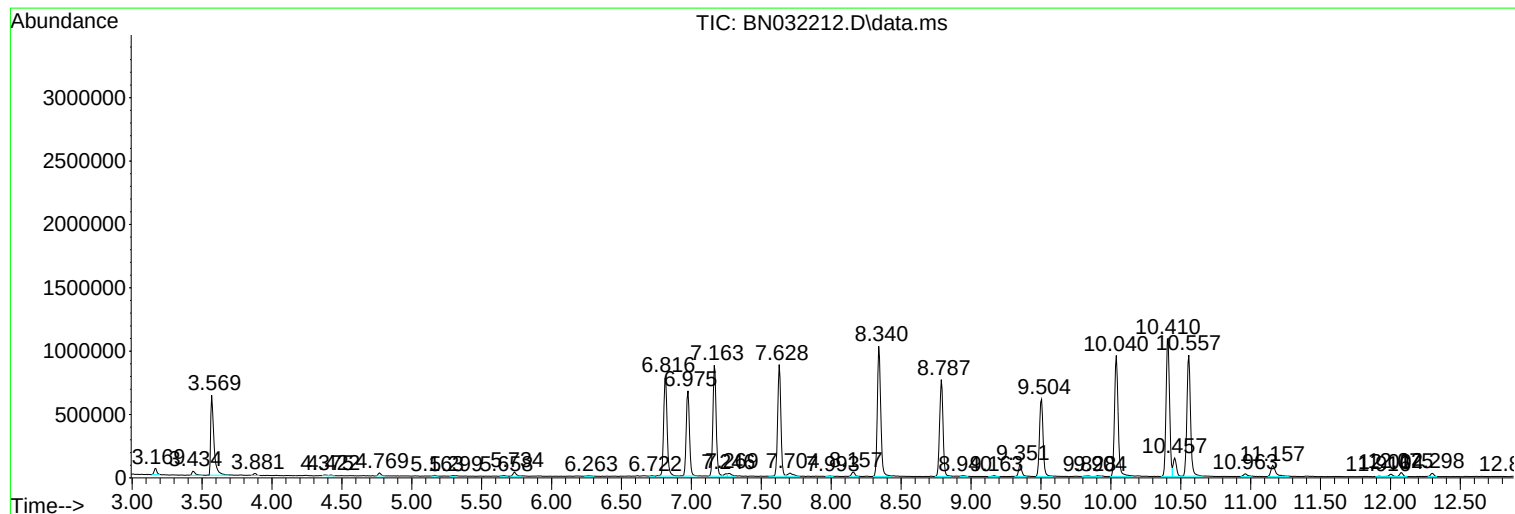
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN052924.MA.M
Quant Title : SVOA CALIBRATION

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



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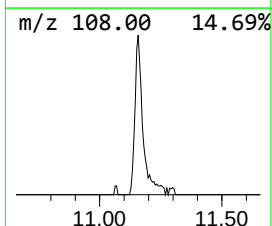
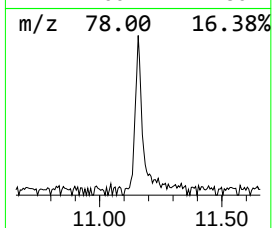
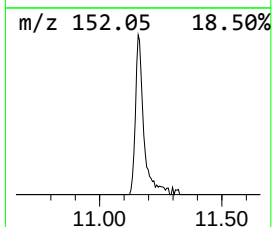
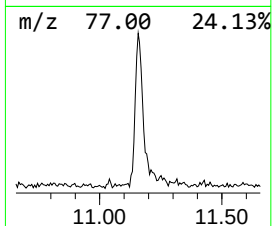
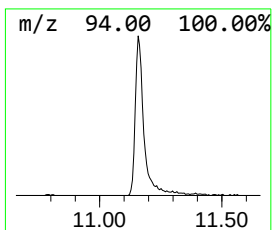
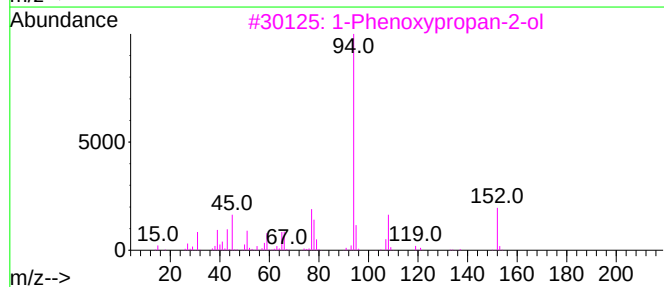
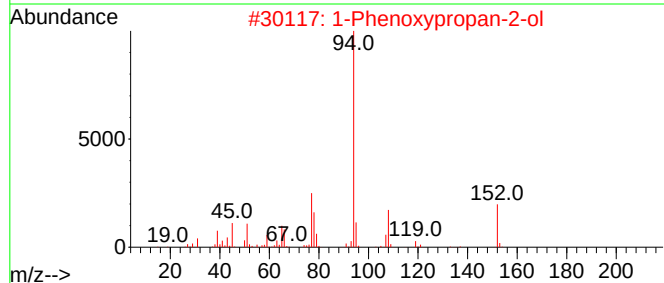
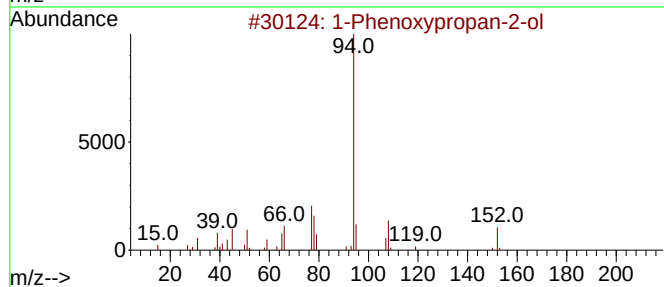
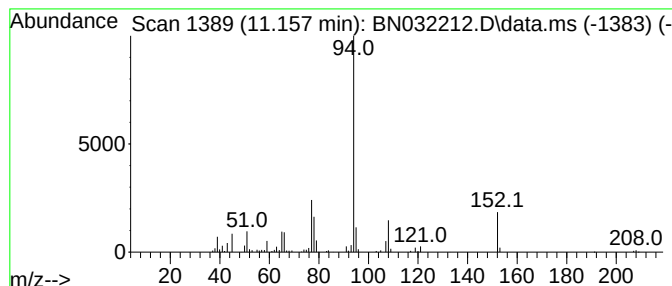
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN052924.MA.M
Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.157	2.24 ng/ul	214809	Naphthalene-d8	10.410

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	97
2			1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	96
3			1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	93
4			1-Propanol, 3-phenoxy-	152	C9H12O2	006180-61-6	91
5			1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	91



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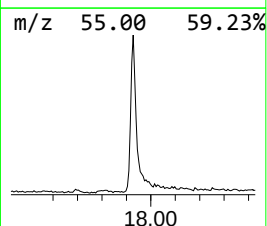
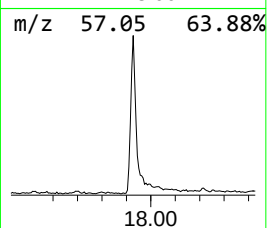
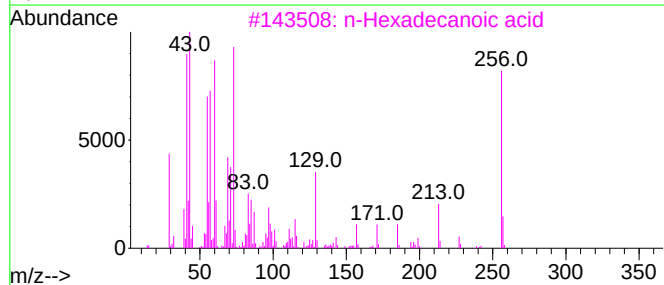
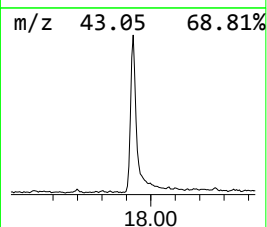
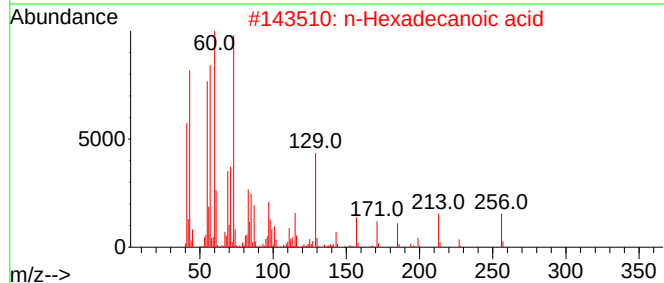
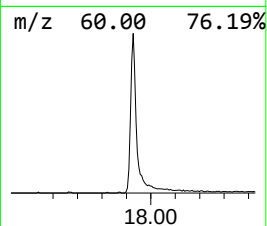
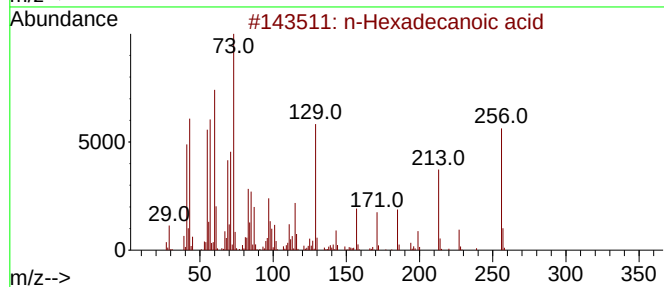
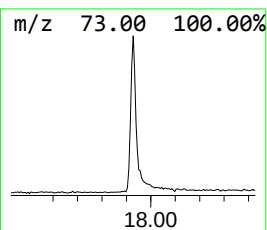
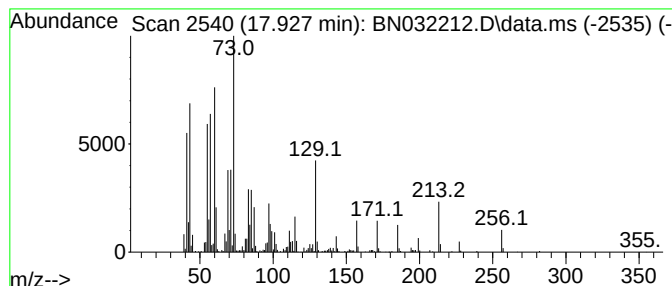
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN052924.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 n-Hexadecanoic acid Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.927	5.99 ng/ul	857153	Phenanthrene-d10	17.028

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			Pentadecanoic acid	242	C15H30O2	001002-84-2	83



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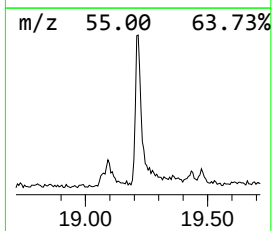
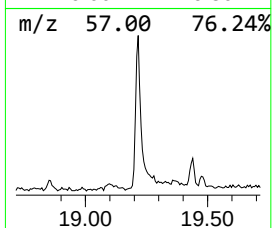
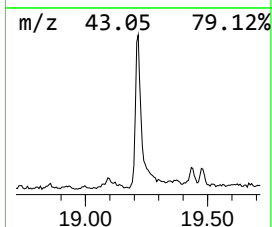
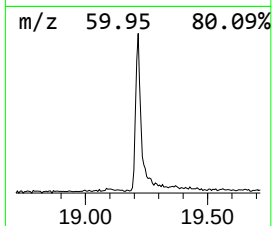
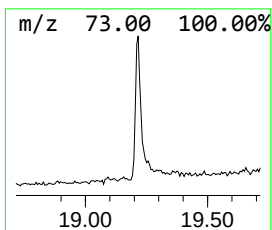
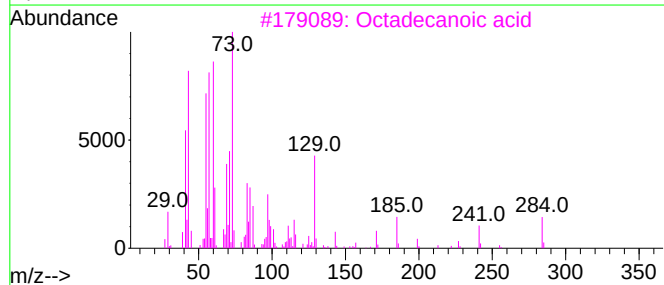
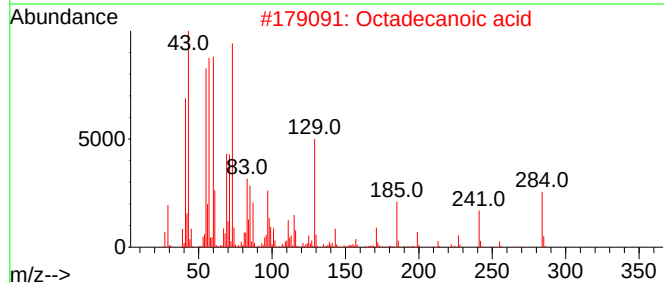
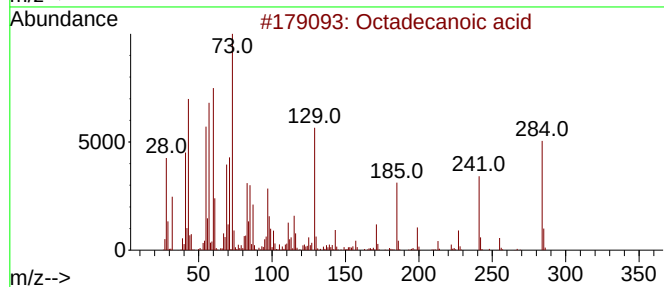
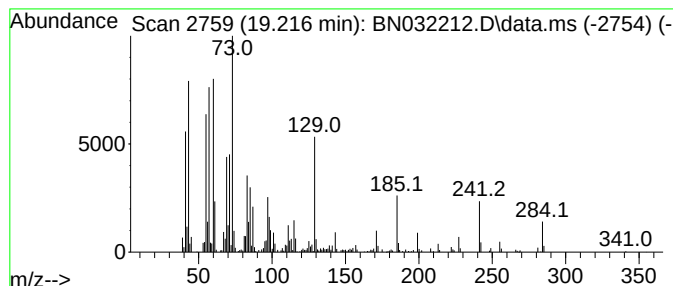
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN052924.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 Octadecanoic acid Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.216	3.20 ng/ul	519308	Chrysene-d12	21.263

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octadecanoic acid	284	C18H36O2	000057-11-4	99
2			Octadecanoic acid	284	C18H36O2	000057-11-4	99
3			Octadecanoic acid	284	C18H36O2	000057-11-4	99
4			Octadecanoic acid	284	C18H36O2	000057-11-4	98
5			Octadecanoic acid	284	C18H36O2	000057-11-4	94



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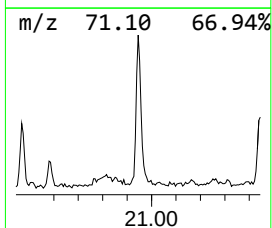
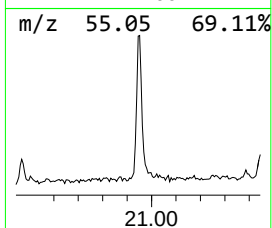
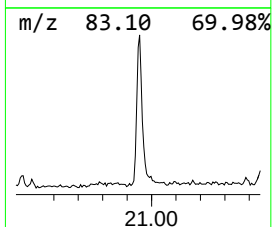
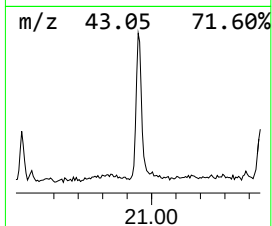
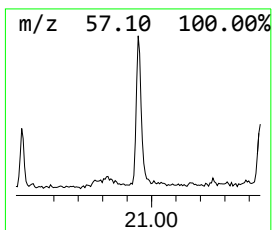
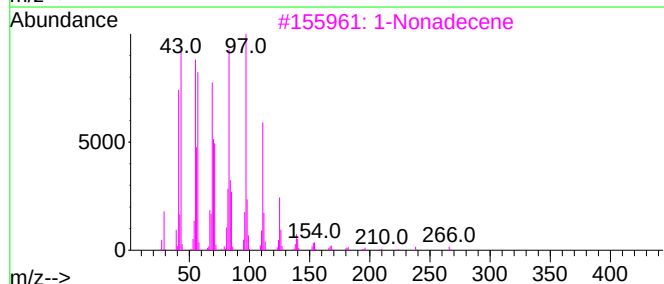
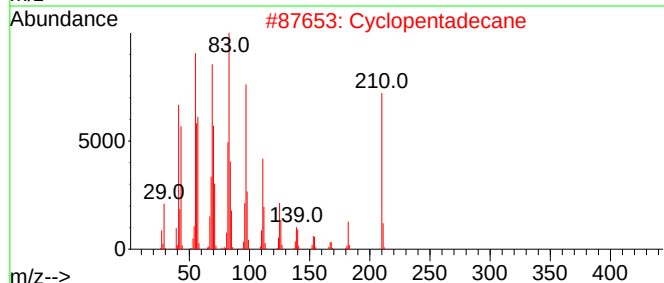
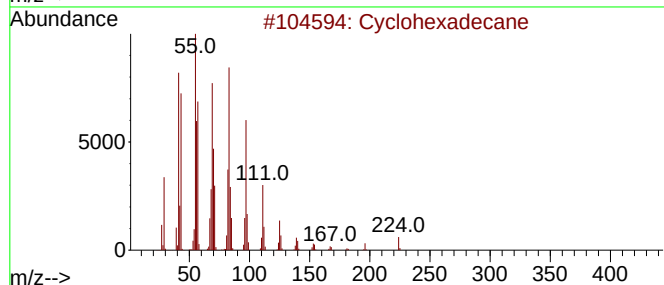
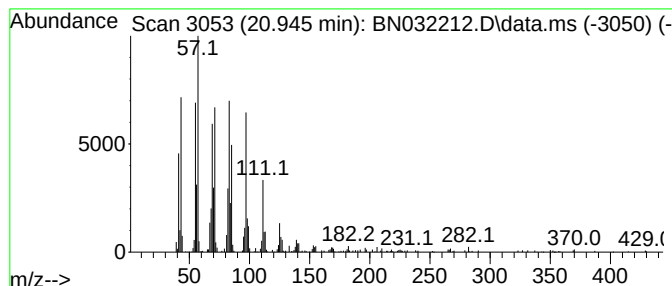
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 (DEL) Alkane: Cyclic20.945 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.945	2.21 ng/ul	358954	Chrysene-d12	21.263

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexadecane	224	C16H32	000295-65-8	98
2			Cyclopentadecane	210	C15H30	000295-48-7	95
3			1-Nonadecene	266	C19H38	018435-45-5	95
4			Pentadecafluorooctanoic acid, oc...	666	C26H37F15O2	1000406-04-8	93
5			1-Heneicosanol	312	C21H44O	015594-90-8	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN062424\
Data File : BN032212.D
Acq On : 24 Jun 2024 17:41
Operator : MA/JU
Sample : P2856-09
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
C0SH1

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
1-Phenoxypropan...	11.157	2.2	ng/u1	214809	2	10.410	1918800	20.0
n-Hexadecanoic ...	17.927	6.0	ng/u1	857153	4	17.028	2860070	20.0
Octadecanoic acid	19.216	3.2	ng/u1	519308	5	21.263	3247000	20.0
(DEL) Alkane: C...	20.945	2.2	ng/u1	358954	5	21.263	3247000	20.0