

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM080924\
 Data File : BM047107.D
 Acq On : 09 Aug 2024 05:31
 Operator : RC/JU
 Sample : P3434-07
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 E05W9

Integration Parameters: LSCINT.P

Integrator: RTE

Smoothing : OFF

Filtering: 5

Sampling : 1

Min Area: 0.1 % of largest Peak

Start Thrs: 0.2

Max Peaks: 100

Stop Thrs : 0

Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >

Peak separation: 5

Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM080224.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BM047107.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.016	18	22	28	rBV	59277	76116	1.55%	0.115%
2	3.275	63	66	69	rBV	108091	133053	2.70%	0.201%
3	3.310	69	72	84	rVV	75595	135386	2.75%	0.204%
4	3.405	84	88	100	rVV	757583	1072904	21.81%	1.617%
5	3.499	100	104	111	rVB	184421	230819	4.69%	0.348%
6	3.704	134	139	145	rVB	27433	38920	0.79%	0.059%
7	4.363	246	251	259	rBV2	20215	39936	0.81%	0.060%
8	4.499	269	274	277	rBV2	21058	30793	0.63%	0.046%
9	4.675	300	304	315	rBV	83922	128314	2.61%	0.193%
10	5.075	364	372	376	rBV2	60935	110820	2.25%	0.167%
11	5.528	444	449	456	rVB2	33227	54797	1.11%	0.083%
12	6.093	540	545	553	rVB2	27856	47187	0.96%	0.071%
13	6.593	623	630	641	rVB	1032464	1604562	32.61%	2.419%
14	6.745	650	656	666	rVB	799096	1191848	24.22%	1.796%
15	6.928	681	687	699	rVB	1034025	1554762	31.60%	2.343%
16	7.028	699	704	711	rBV	52966	80647	1.64%	0.122%
17	7.387	756	765	772	rBV	918722	1407259	28.60%	2.121%
18	7.481	776	781	788	rVB5	18882	37651	0.77%	0.057%
19	7.892	846	851	857	rBV5	21927	36104	0.73%	0.054%
20	8.010	862	871	877	rBV	231863	423427	8.61%	0.638%
21	8.104	881	887	896	rBV	1184428	1847537	37.55%	2.785%
22	8.481	946	951	954	rBV2	37308	56567	1.15%	0.085%
23	8.528	954	959	967	rVV	949208	1524351	30.98%	2.298%
24	8.610	967	973	981	rVV	2129108	3030476	61.60%	4.568%
25	9.116	1049	1059	1065	rBV	119948	194555	3.95%	0.293%
26	9.239	1074	1080	1087	rBV	812344	1291792	26.26%	1.947%
27	9.428	1107	1112	1119	rBV2	28688	61503	1.25%	0.093%
28	9.781	1161	1172	1182	rBV	1245510	2074998	42.18%	3.128%
29	10.139	1224	1233	1240	rBV	1219848	1967626	39.99%	2.966%
30	10.292	1248	1259	1265	rBV	748589	1282955	26.08%	1.934%
31	10.345	1265	1268	1283	rVB	56758	123412	2.51%	0.186%
32	10.716	1325	1331	1336	rBV2	44025	69869	1.42%	0.105%
33	10.898	1355	1362	1372	rBV	184086	346515	7.04%	0.522%
34	11.316	1427	1433	1439	rBV2	80925	128288	2.61%	0.193%
35	11.598	1475	1481	1486	rBV2	29292	45827	0.93%	0.069%
36	11.675	1489	1494	1499	rVV2	24604	41883	0.85%	0.063%

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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_M\Methods\SFAM-EPA-BM080224.MA.M
 Title : SVOA CALIBRATION

37	11.739	1499	1505	1512	rVB2	27470	47386	0.96%	0.071%
38	12.492	1627	1633	1636	rBV3	29557	53862	1.09%	0.081%
39	12.527	1636	1639	1644	rVV5	37429	55499	1.13%	0.084%
40	12.757	1673	1678	1682	rBV7	27287	45156	0.92%	0.068%
41	12.939	1701	1709	1713	rVV4	24840	49756	1.01%	0.075%
42	12.986	1713	1717	1719	rVV	17157	25322	0.51%	0.038%
43	13.016	1719	1722	1727	rVB	38012	51684	1.05%	0.078%
44	13.074	1727	1732	1737	rBV6	16584	29692	0.60%	0.045%
45	13.151	1740	1745	1754	rBV5	63351	118436	2.41%	0.179%
46	13.469	1792	1799	1804	rVV	2026771	2784162	56.59%	4.197%
47	13.504	1804	1805	1809	rVV	96694	76672	1.56%	0.116%
48	13.727	1836	1843	1854	rBV	2324164	3424368	69.60%	5.162%
49	13.892	1863	1871	1878	rVB3	46619	75118	1.53%	0.113%
50	14.039	1888	1896	1904	rBV	1823745	2627283	53.40%	3.960%
51	14.286	1931	1938	1952	rBV	845732	1468957	29.86%	2.214%
52	14.427	1957	1962	1970	rVV3	51285	94142	1.91%	0.142%
53	14.874	2035	2038	2045	rBV4	18664	29348	0.60%	0.044%
54	15.045	2060	2067	2074	rBV	3011787	4206693	85.50%	6.341%
55	15.127	2077	2081	2085	rVB4	25288	32913	0.67%	0.050%
56	15.186	2085	2091	2099	rBV	648497	913353	18.56%	1.377%
57	15.286	2104	2108	2113	rVB7	16963	30925	0.63%	0.047%
58	15.539	2142	2151	2159	rBV7	19601	45868	0.93%	0.069%
59	15.639	2161	2168	2171	rBV8	22278	45161	0.92%	0.068%
60	15.745	2181	2186	2190	rBV5	24463	38521	0.78%	0.058%
61	15.951	2218	2221	2229	rVB9	15171	31521	0.64%	0.048%
62	16.245	2263	2271	2278	rBV9	22653	50918	1.03%	0.077%
63	16.598	2327	2331	2334	rVV5	20875	31924	0.65%	0.048%
64	16.627	2334	2336	2342	rVV7	16352	25017	0.51%	0.038%
65	16.686	2342	2346	2351	rVB2	27198	37090	0.75%	0.056%
66	16.739	2351	2355	2359	rBV3	61066	88385	1.80%	0.133%
67	16.798	2360	2365	2371	rVV	2116605	3024188	61.47%	4.558%
68	16.839	2371	2372	2376	rVV	50405	51515	1.05%	0.078%
69	16.898	2376	2382	2394	rVB2	2374571	3229969	65.65%	4.868%
70	17.009	2396	2401	2406	rVB3	22874	45375	0.92%	0.068%
71	17.104	2414	2417	2422	rVV6	16979	24873	0.51%	0.037%
72	17.221	2434	2437	2442	rVV6	17287	25359	0.52%	0.038%
73	17.486	2477	2482	2485	rBV	103814	142929	2.91%	0.215%
74	17.574	2492	2497	2501	rBV3	36677	68977	1.40%	0.104%
75	17.615	2501	2504	2509	rVB	62416	82712	1.68%	0.125%
76	17.674	2509	2514	2520	rBV	234792	332198	6.75%	0.501%

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77	17.739	2520	2525	2533	rBV	944810	1349805	27.44%	2.035%
78	17.992	2564	2568	2571	rBV3	37142	52589	1.07%	0.079%
79	18.033	2571	2575	2583	rVV2	183732	293575	5.97%	0.443%
80	18.192	2596	2602	2607	rBV2	226394	331204	6.73%	0.499%
81	18.433	2638	2643	2647	rBV	81494	118903	2.42%	0.179%
82	18.515	2653	2657	2660	rVB	70896	82912	1.69%	0.125%
83	18.609	2671	2673	2679	rVB7	19897	25433	0.52%	0.038%
84	18.768	2697	2700	2707	rVB	50190	72328	1.47%	0.109%
85	18.845	2709	2713	2715	rBV2	64829	90212	1.83%	0.136%
86	18.880	2715	2719	2722	rBV	79883	102645	2.09%	0.155%
87	18.915	2722	2725	2727	rBV	86877	96580	1.96%	0.146%
88	19.039	2742	2746	2754	rBV	180655	277067	5.63%	0.418%
89	19.215	2769	2776	2788	rBV	3484548	4919928	100.00%	7.416%
90	19.303	2789	2791	2798	rVB6	35704	42784	0.87%	0.064%
91	19.433	2807	2813	2821	rBV	1928445	2384958	48.48%	3.595%
92	19.598	2838	2841	2843	rBV3	44022	41866	0.85%	0.063%
93	19.768	2867	2870	2874	rBV2	88599	112352	2.28%	0.169%
94	20.174	2936	2939	2943	rVB	112901	132633	2.70%	0.200%
95	20.262	2951	2954	2957	rBV2	75183	84446	1.72%	0.127%
96	20.327	2963	2965	2970	rVB2	75978	95280	1.94%	0.144%
97	20.774	3037	3041	3052	rVB	558491	777639	15.81%	1.172%
98	21.033	3080	3085	3090	rBV2	2241274	3007352	61.13%	4.533%
99	23.550	3506	3513	3522	rVB	1433057	3090845	62.82%	4.659%
100	23.733	3537	3544	3552	rBV	1173780	2574349	52.32%	3.880%

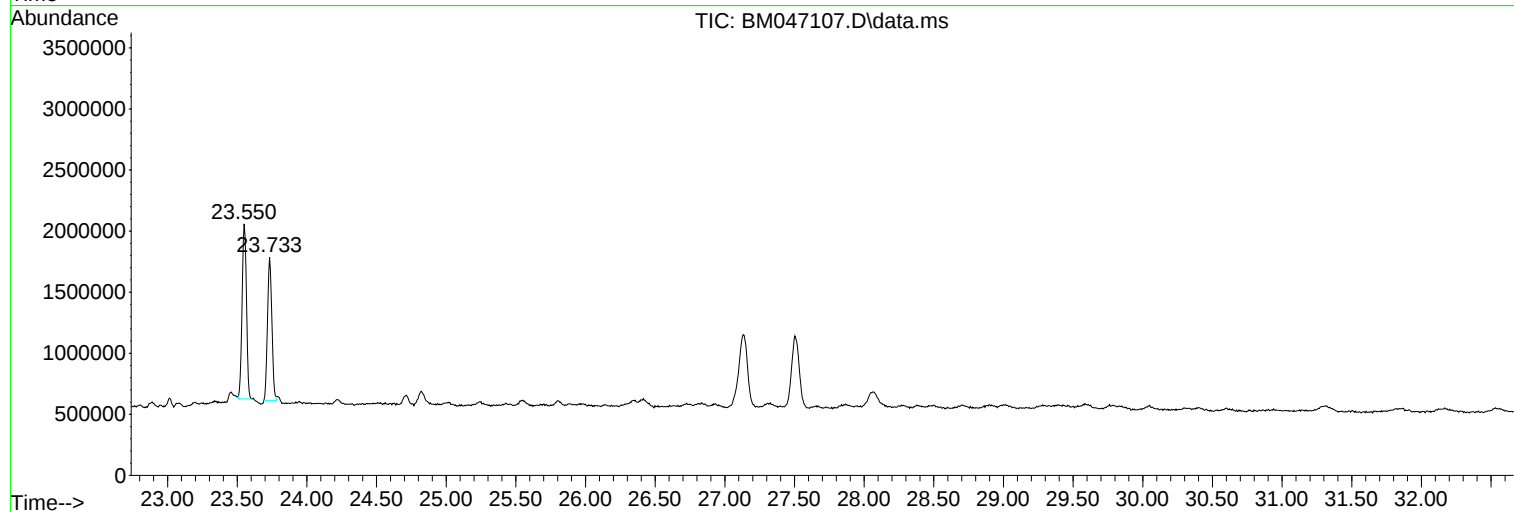
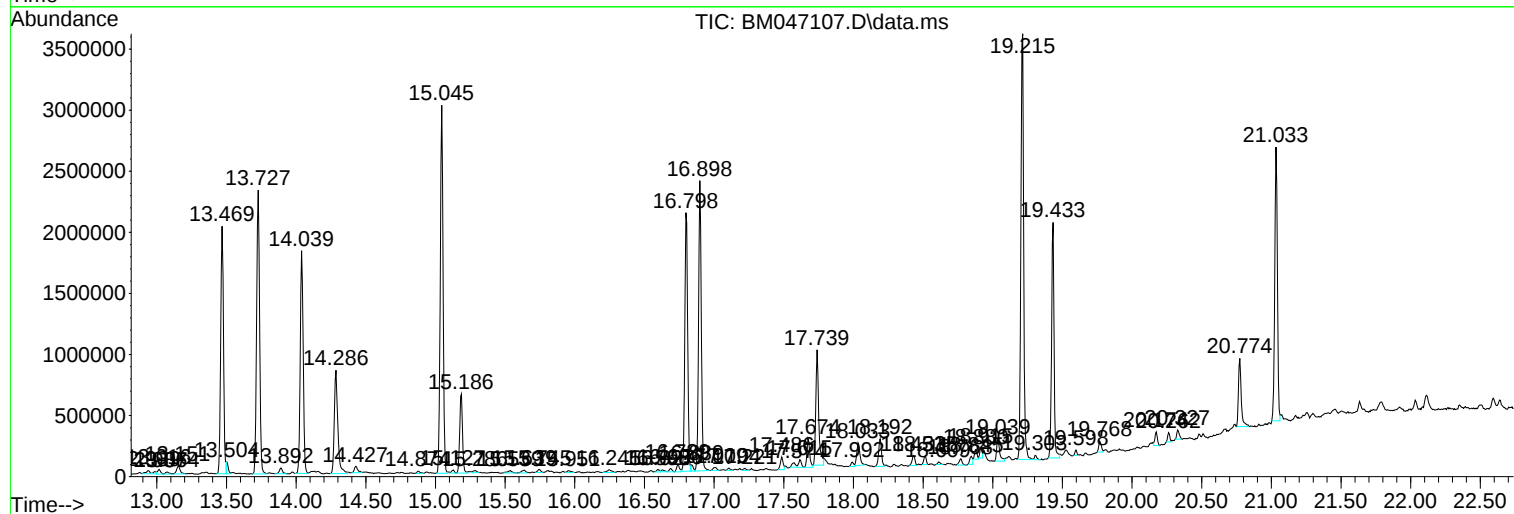
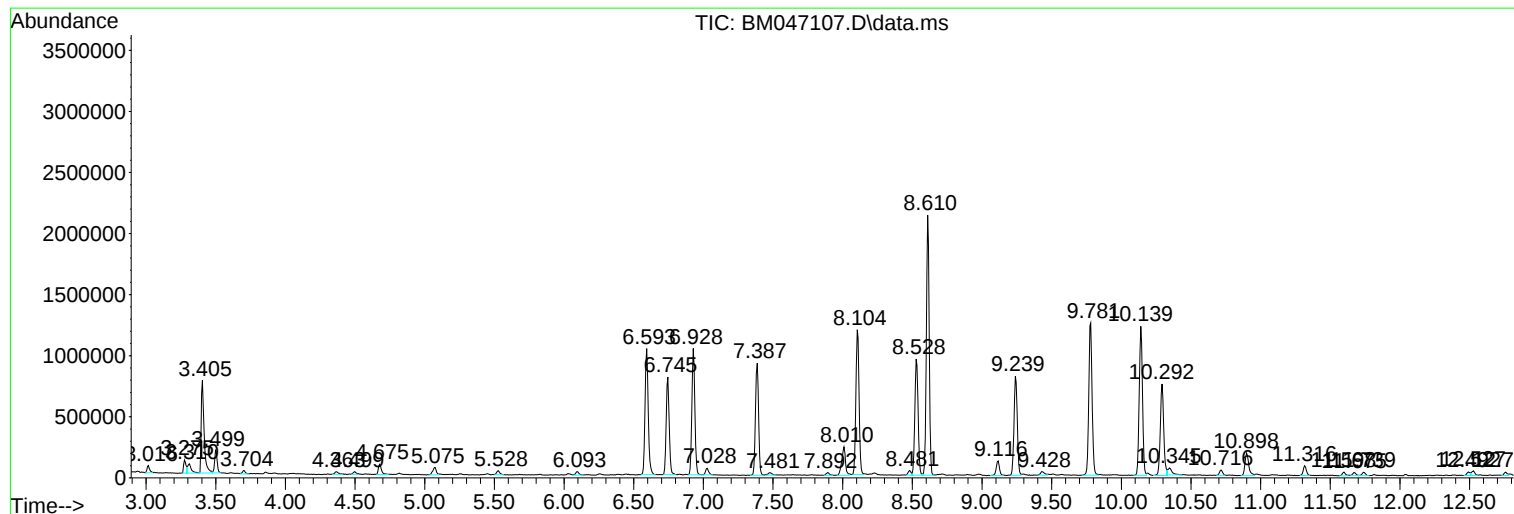
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM080224.MA.M
Quant Title : SVOA CALIBRATION

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



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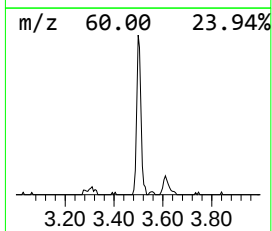
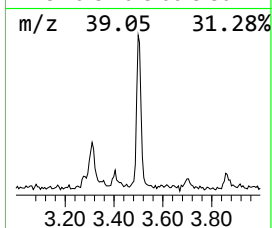
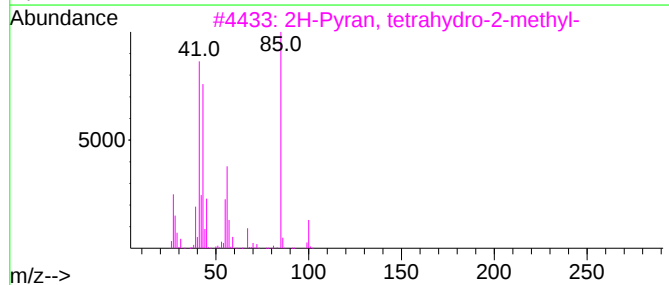
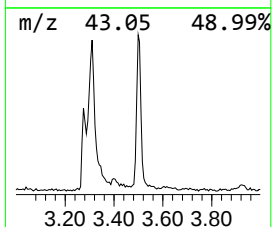
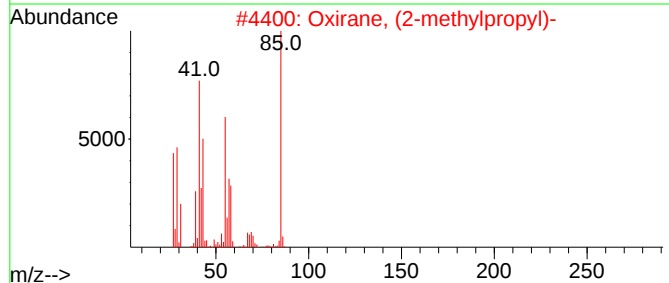
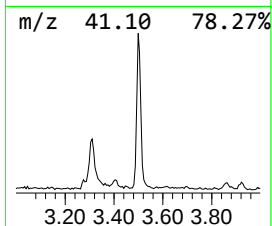
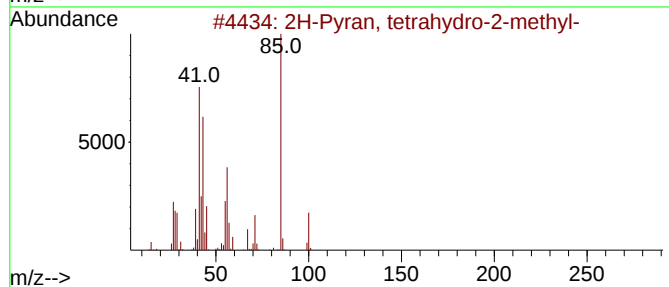
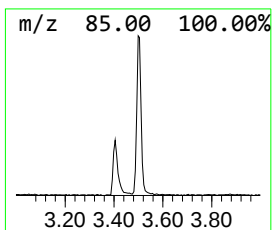
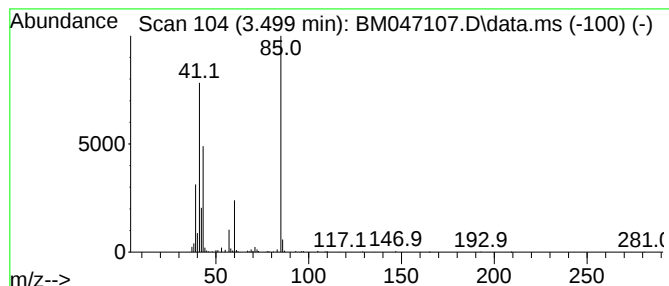
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM080224.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 unknown-01 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.499	3.28 ng/ul	230819	1,4-Dichlorobenzene-d4	7.387

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2H-Pyran, tetrahydro-2-methyl-	100	C6H12O	010141-72-7	38	
2	Oxirane, (2-methylpropyl)-	100	C6H12O	023850-78-4	36	
3	2H-Pyran, tetrahydro-2-methyl-	100	C6H12O	010141-72-7	28	
4	2-Pyrrolidinone	85	C4H7NO	000616-45-5	28	
5	5-Hydroxypentanal	102	C5H10O2	004221-03-8	25	



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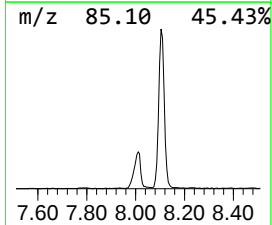
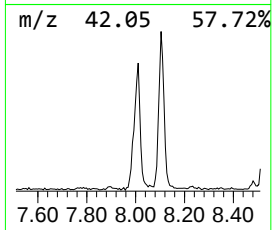
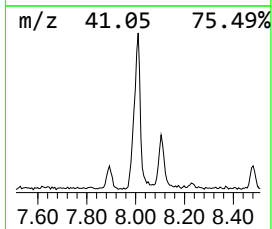
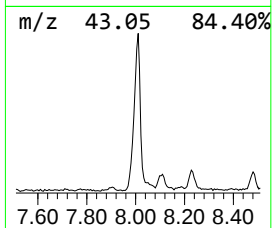
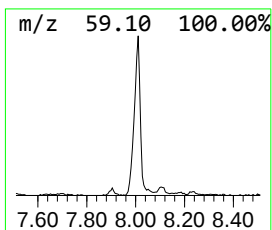
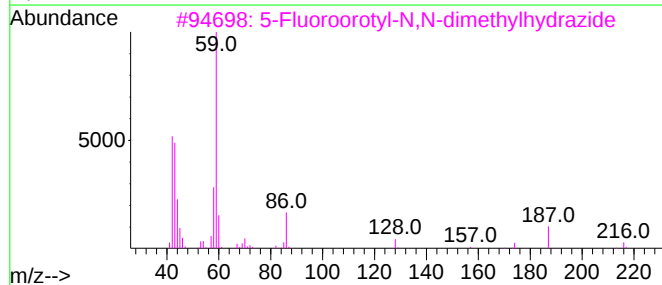
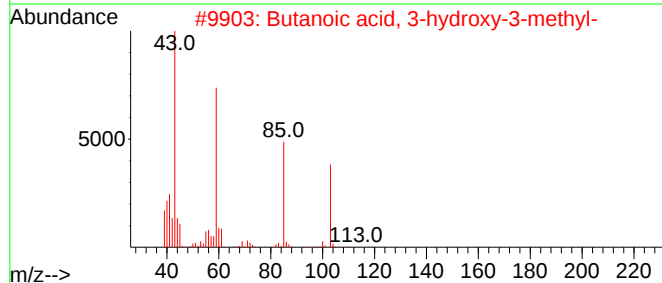
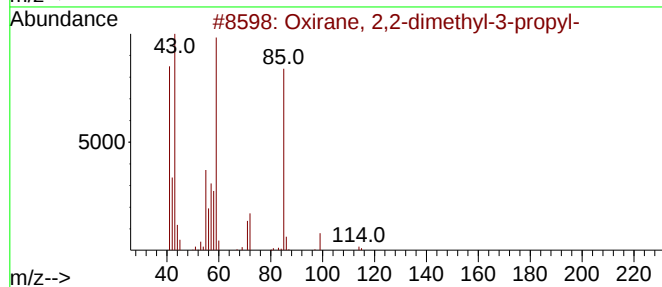
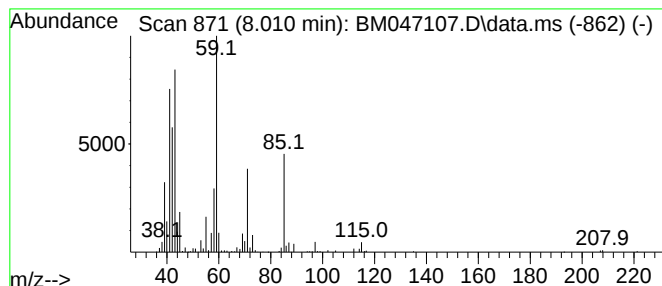
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM080224.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 Oxirane, 2,2-dimethyl-3-pro... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.010	6.02 ng/ul	423427	1,4-Dichlorobenzene-d4	7.387

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Oxirane, 2,2-dimethyl-3-propyl-	114	C7H14O	017612-35-0	53
2			Butanoic acid, 3-hydroxy-3-methyl-	118	C5H10O3	000625-08-1	38
3			5-Fluoroorotyl-N,N-dimethylhydra...	216	C7H9FN4O3	1000227-24-4	37
4			Tetraglyme	222	C10H22O5	000143-24-8	35
5			Tetraglyme	222	C10H22O5	000143-24-8	35



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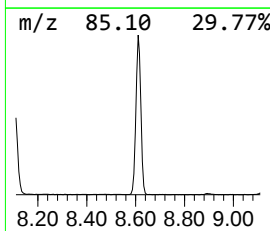
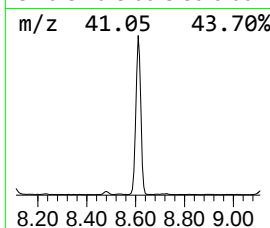
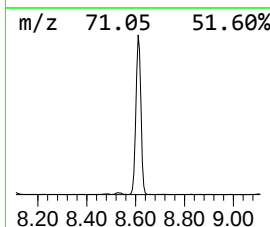
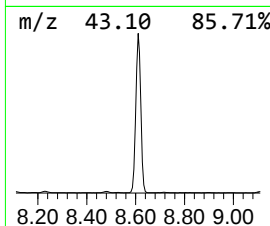
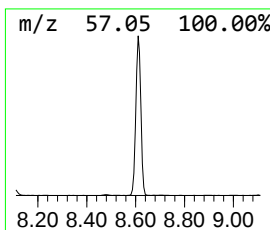
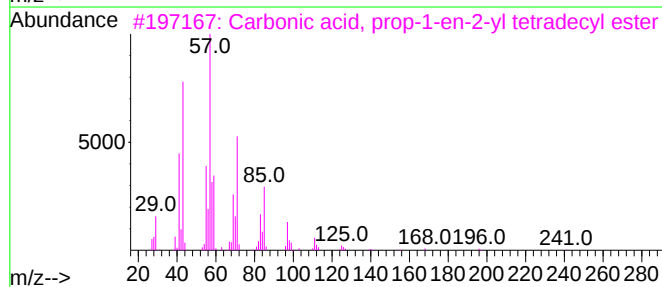
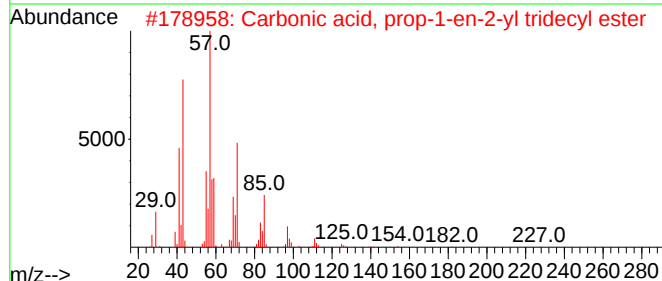
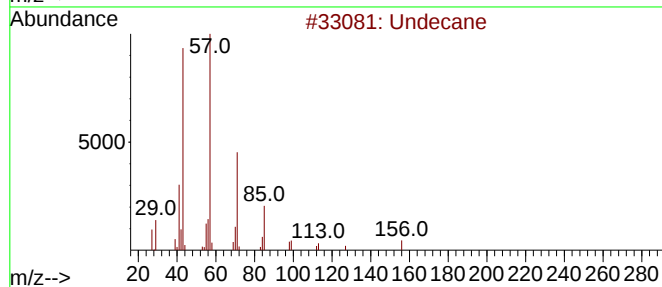
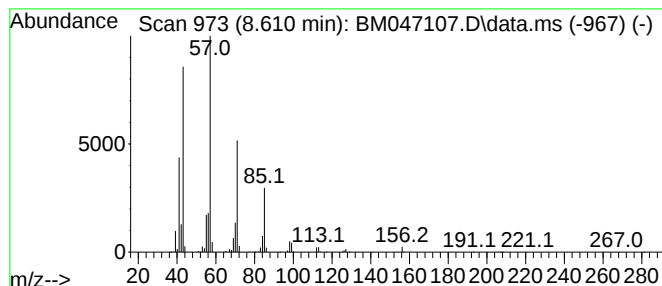
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM080224.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 (DEL) Alkane: Straight-Chai... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.610	43.07 ng/ul	3030480	1,4-Dichlorobenzene-d4	7.387

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Undecane	156	C11H24	001120-21-4	91
2			Carbonic acid, prop-1-en-2-yl tr...	284	C17H32O3	1000382-90-8	90
3			Carbonic acid, prop-1-en-2-yl te...	298	C18H34O3	1000382-90-9	90
4			Tetradecane	198	C14H30	000629-59-4	90
5			Hexadecane	226	C16H34	000544-76-3	78



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM080924\
Data File : BM047107.D
Acq On : 09 Aug 2024 05:31
Operator : RC/JU
Sample : P3434-07
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
E05W9

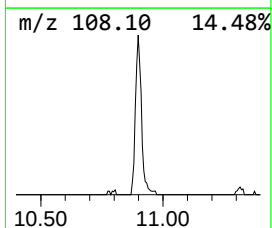
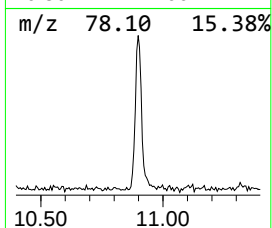
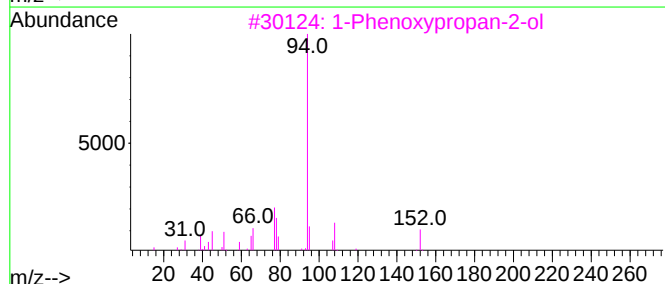
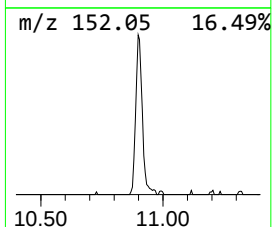
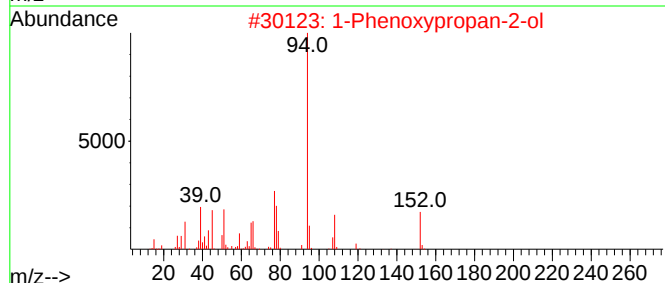
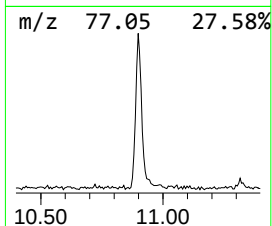
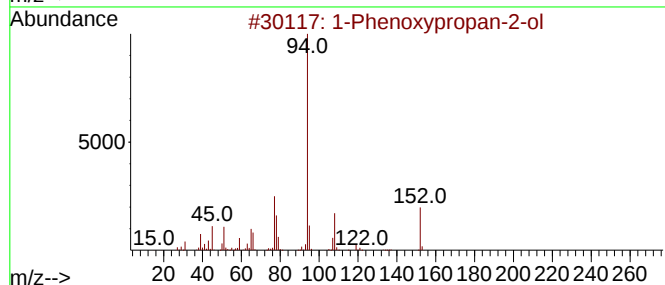
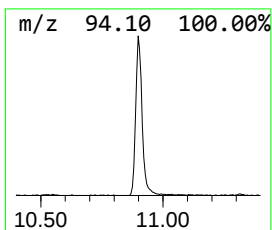
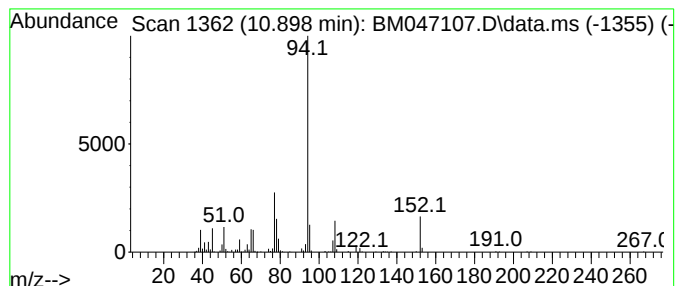
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 1-Phenoxypropan-2-ol Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.898	3.52 ng/ul	346515	Naphthalene-d8	10.139

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-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	91
4			1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	87
5			1-Propanol, 3-phenoxy-	152	C9H12O2	006180-61-6	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM080924\
Data File : BM047107.D
Acq On : 09 Aug 2024 05:31
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Sample : P3434-07
Misc :
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ClientSampleId :
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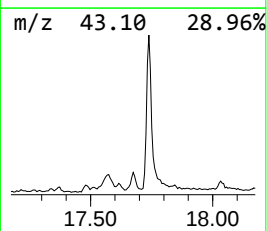
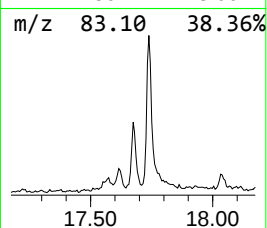
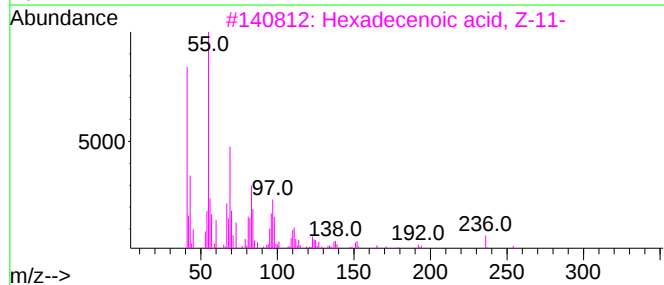
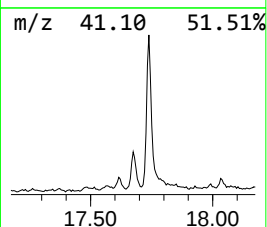
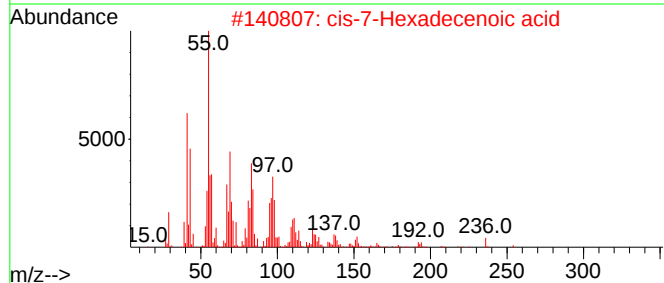
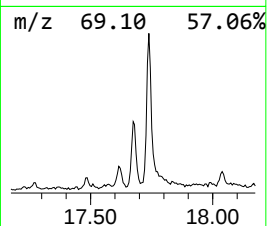
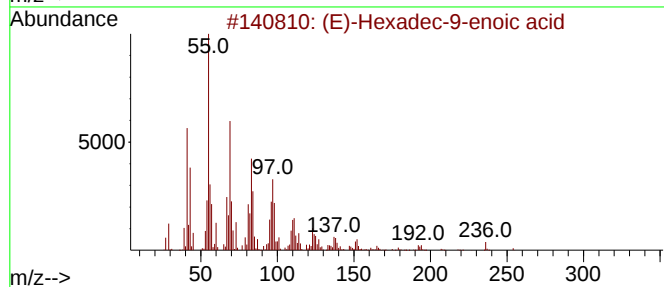
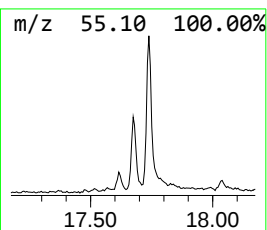
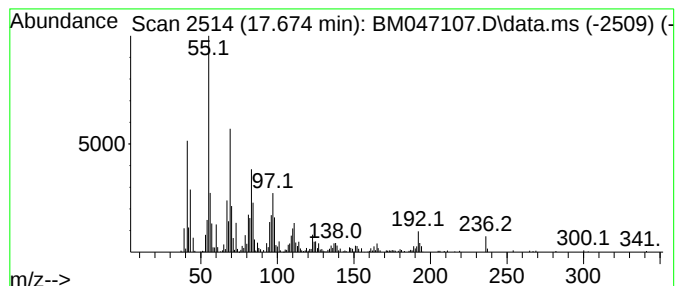
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 (E)-Hexadec-9-enoic acid Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.674	2.20 ng/ul	332198	Phenanthrene-d10	16.798

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			(E)-Hexadec-9-enoic acid	254	C16H30O2	010030-73-6	98
2			cis-7-Hexadecenoic acid	254	C16H30O2	002416-19-5	98
3			Hexadecenoic acid, Z-11-	254	C16H30O2	002416-20-8	91
4			Palmitoleic acid	254	C16H30O2	000373-49-9	90
5			Oxacycloheptadecan-2-one	254	C16H30O2	000109-29-5	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM080924\
Data File : BM047107.D
Acq On : 09 Aug 2024 05:31
Operator : RC/JU
Sample : P3434-07
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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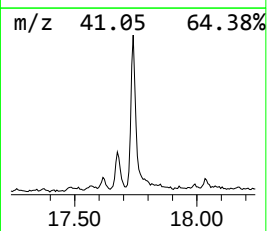
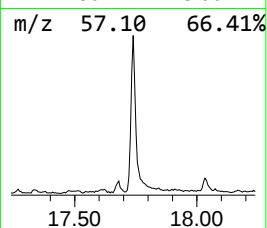
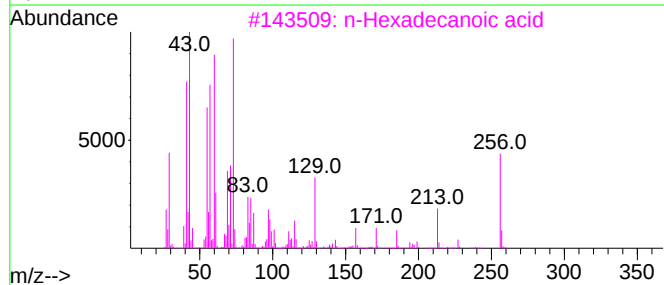
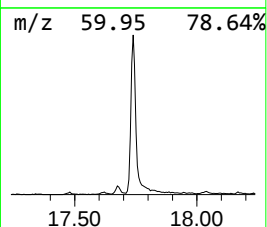
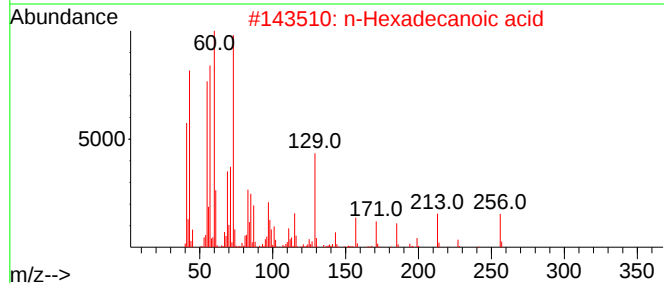
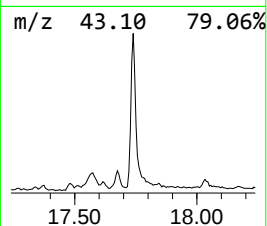
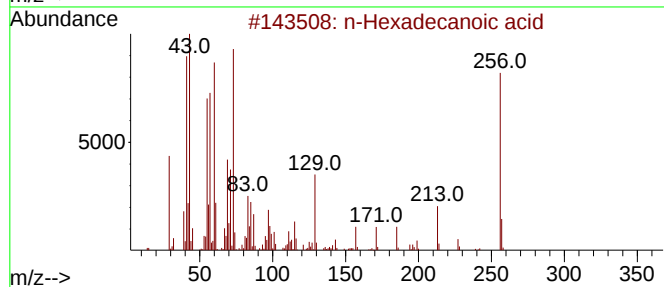
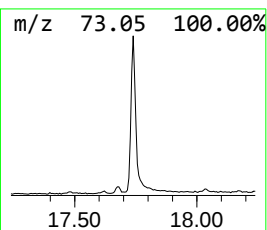
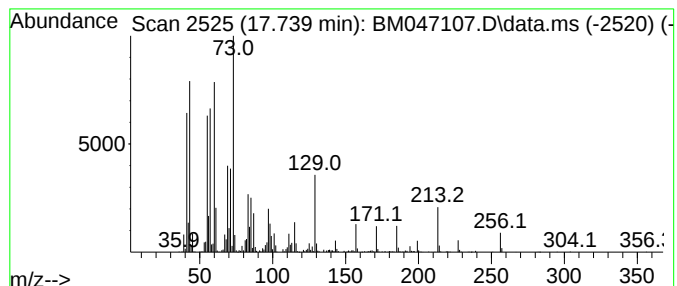
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 n-Hexadecanoic acid Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.739	8.93 ng/ul	1349810	Phenanthrene-d10	16.798

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			Tetradecanoic acid	228	C14H28O2	000544-63-8	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM080924\
Data File : BM047107.D
Acq On : 09 Aug 2024 05:31
Operator : RC/JU
Sample : P3434-07
Misc :
ALS Vial : 24 Sample Multiplier: 1

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ClientSampleId :
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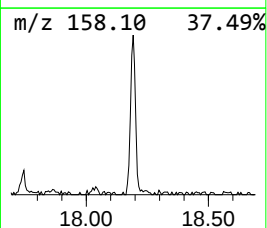
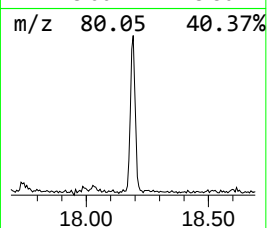
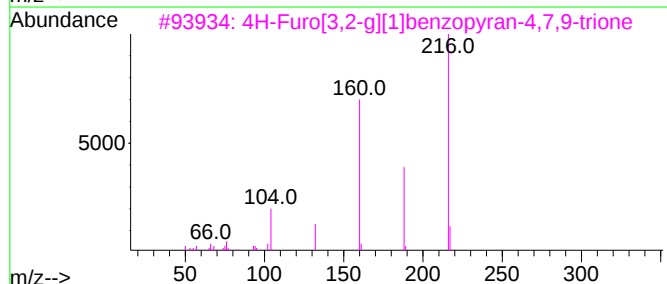
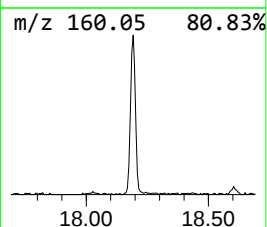
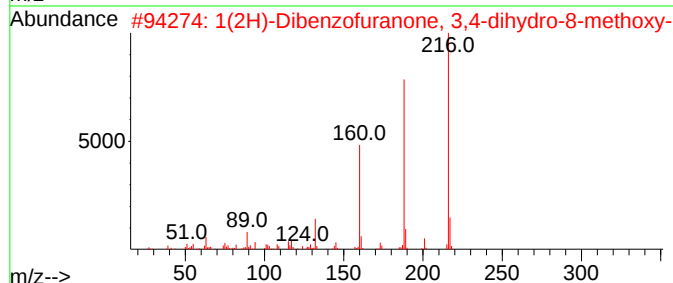
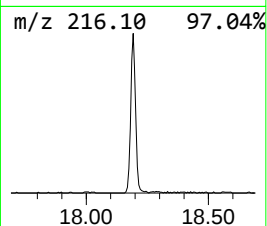
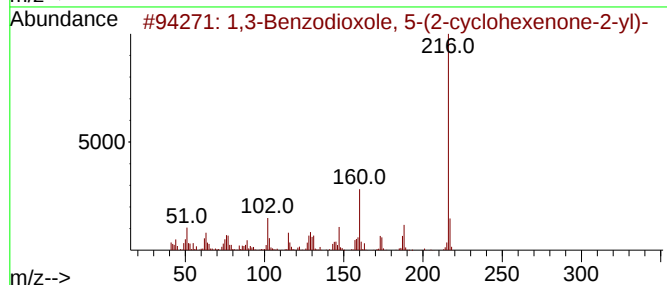
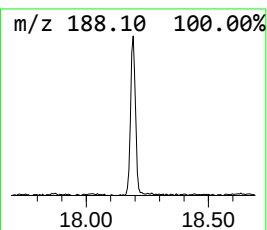
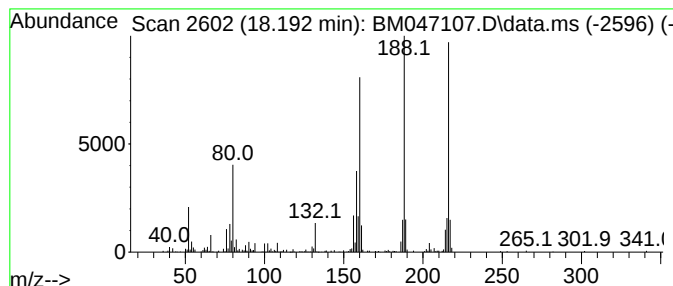
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 unknown-02 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.192	2.19 ng/ul	331204	Phenanthrene-d10	16.798

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1,3-Benzodioxole, 5-(2-cyclohexe...	216	C13H12O3	042866-84-2	43
2			1(2H)-Dibenzofuranone, 3,4-dihyd...	216	C13H12O3	095338-40-2	43
3			4H-Furo[3,2-g][1]benzopyran-4,7,...	216	C11H4O5	000483-36-3	38
4			4-(4-Ethoxyphenyl)-1H-imidazole	188	C11H12N2O	1000311-54-8	25
5			1,2-Dihydro-3-isopropyl-2-oxoqui...	188	C11H12N2O	015054-64-5	22



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM080924\
Data File : BM047107.D
Acq On : 09 Aug 2024 05:31
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Sample : P3434-07
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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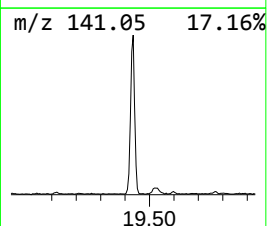
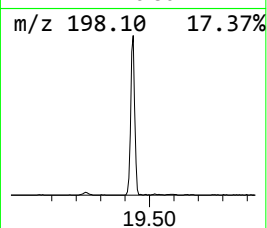
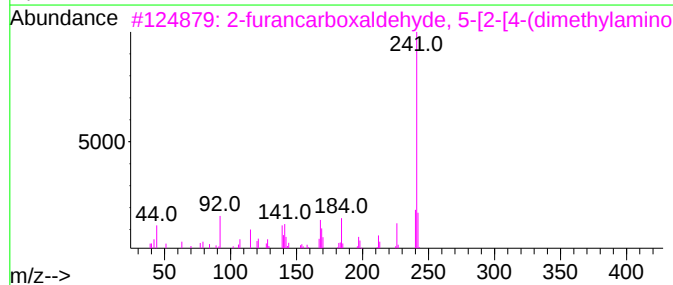
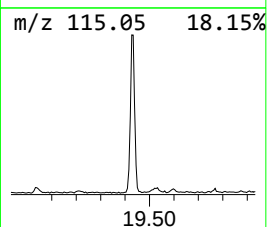
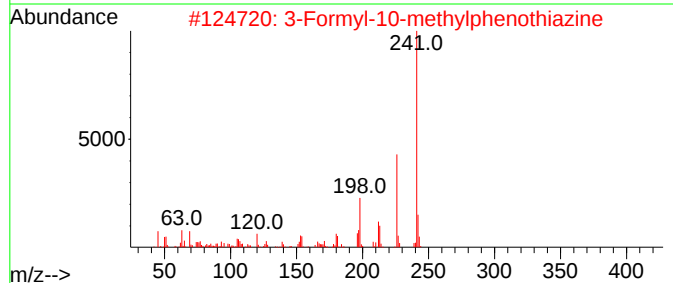
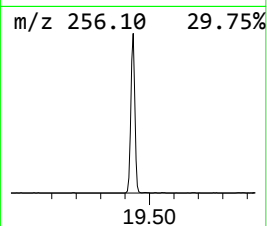
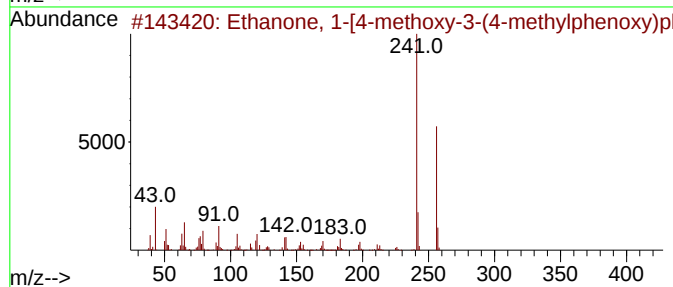
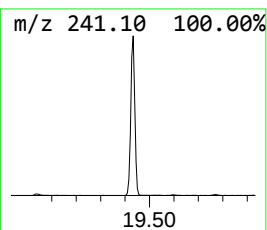
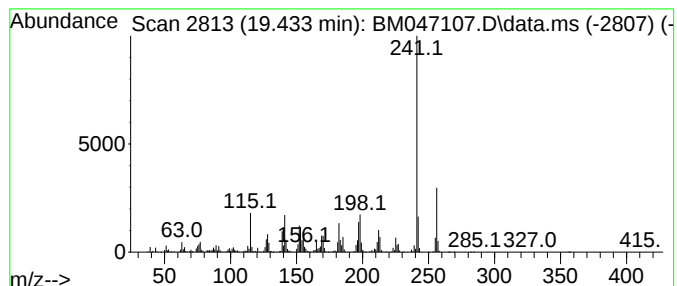
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 Ethanone, 1-[4-methoxy-3-(4... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.433	15.86 ng/ul	2384960	Chrysene-d12	21.033

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Ethanone, 1-[4-methoxy-3-(4-meth...	256	C16H16O3	116345-94-9	62
2			3-Formyl-10-methylphenothiazine	241	C14H11NOS	004997-36-8	58
3			2-furancarboxaldehyde, 5-[2-[4-(...	241	C15H15NO2	1000396-52-3	58
4			4-Isopropylamino-N-isopropylbenz...	256	C12H20N2O2S	1010395-71-0	53
5			Bisphenol C	256	C17H20O2	000079-97-0	52



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM080924\
Data File : BM047107.D
Acq On : 09 Aug 2024 05:31
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Sample : P3434-07
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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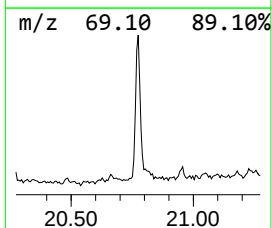
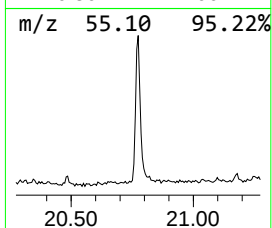
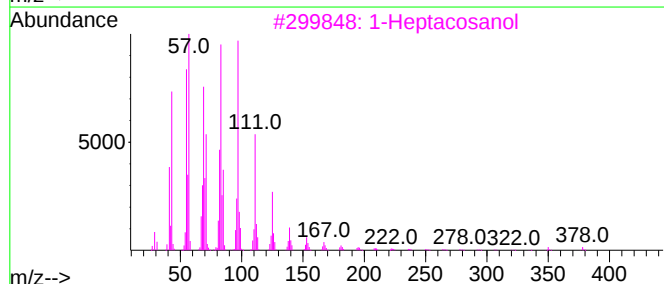
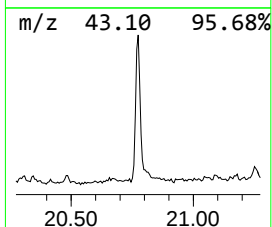
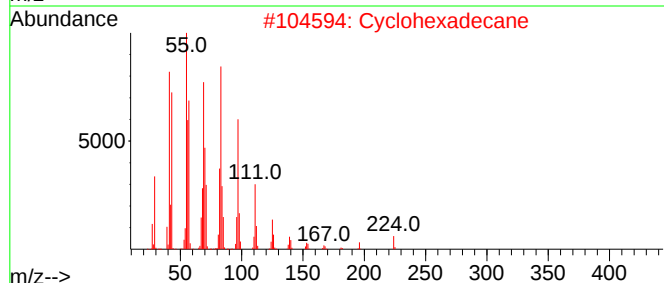
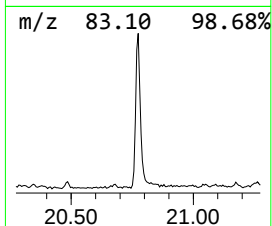
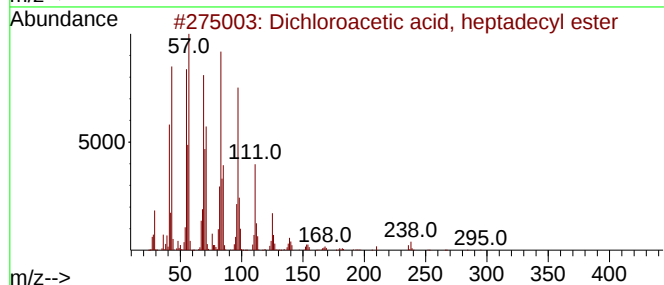
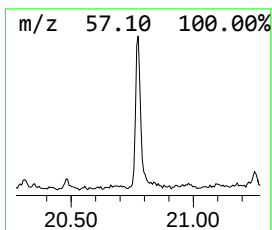
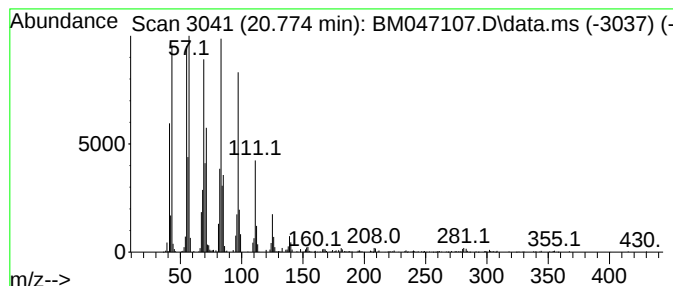
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 Dichloroacetic acid, heptad... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.774	5.17 ng/ul	777639	Chrysene-d12	21.033

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Dichloroacetic acid, heptadecyl ...	366	C19H36Cl2O2	1000282-98-2	95
2			Cyclohexadecane	224	C16H32	000295-65-8	94
3			1-Heptacosanol	396	C27H56O	002004-39-9	94
4			1-Heneicosyl formate	340	C22H44O2	077899-03-7	94
5			Behenic alcohol	326	C22H46O	000661-19-8	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM080924\
Data File : BM047107.D
Acq On : 09 Aug 2024 05:31
Operator : RC/JU
Sample : P3434-07
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM080224.MA.M
Quant Title : SVOA CALIBRATION

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

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					#	RT	Resp	Conc
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Oxirane, 2,2-di...	8.010	6.0	ng/ul	423427	1	7.387	1407260	20.0
(DEL) Alkane: S...	8.610	43.1	ng/ul	3030480	1	7.387	1407260	20.0
1-Phenoxypropan...	10.898	3.5	ng/ul	346515	2	10.139	1967630	20.0
(E)-Hexadec-9-e...	17.674	2.2	ng/ul	332198	4	16.798	3024190	20.0
n-Hexadecanoic ...	17.739	8.9	ng/ul	1349810	4	16.798	3024190	20.0
unknown-02	18.192	2.2	ng/ul	331204	4	16.798	3024190	20.0
Ethanone, 1-[4-...	19.433	15.9	ng/ul	2384960	5	21.033	3007350	20.0
Dichloroacetic ...	20.774	5.2	ng/ul	777639	5	21.033	3007350	20.0