

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP013125\
 Data File : BP023844.D
 Acq On : 01 Feb 2025 04:16
 Operator : RC/JU
 Sample : Q1224-08
 Misc :
 ALS Vial : 28 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 A6306

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_P\Methods\SFAM-EPA-BP012925.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BP023844.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.270	27	31	33	rBV	34399	46779	0.69%	0.048%
2	3.293	33	35	40	rVB	63372	75965	1.13%	0.078%
3	3.711	103	106	115	rBV5	16447	26452	0.39%	0.027%
4	3.805	118	122	127	rVB	90187	113882	1.69%	0.116%
5	4.028	156	160	166	rVB	23729	34818	0.52%	0.036%
6	4.122	172	176	183	rVB2	40295	52671	0.78%	0.054%
7	4.393	218	222	227	rVB6	14201	23140	0.34%	0.024%
8	4.552	244	249	259	rVB3	43185	78961	1.17%	0.081%
9	4.934	308	314	321	rBV2	109033	173331	2.57%	0.177%
10	6.187	522	527	533	rBV2	32352	46576	0.69%	0.048%
11	6.793	624	630	636	rBV3	16915	29728	0.44%	0.030%
12	6.964	648	659	675	rVB	1127718	1996412	29.64%	2.041%
13	7.116	678	685	696	rBV	883278	1375118	20.42%	1.406%
14	7.322	713	720	731	rVV	1158129	1786087	26.52%	1.826%
15	7.781	791	798	805	rVV	1453447	2300014	34.15%	2.351%
16	7.846	805	809	815	rVB2	24957	42497	0.63%	0.043%
17	8.487	908	918	926	rBV	750814	1178360	17.50%	1.204%
18	8.587	929	935	945	rVB	352560	568145	8.44%	0.581%
19	8.922	985	992	1003	rVV	986605	1639637	24.34%	1.676%
20	9.363	1062	1067	1074	rVB	37903	60951	0.90%	0.062%
21	9.481	1081	1087	1094	rBV	152724	222785	3.31%	0.228%
22	9.640	1106	1114	1120	rBV	791318	1330658	19.76%	1.360%
23	9.693	1120	1123	1128	rVV2	44889	69719	1.04%	0.071%
24	9.752	1128	1133	1138	rVV	40044	70557	1.05%	0.072%
25	9.805	1138	1142	1150	rVV	67905	127016	1.89%	0.130%
26	9.881	1150	1155	1162	rVB5	11840	22831	0.34%	0.023%
27	9.999	1167	1175	1181	rVB2	25175	51018	0.76%	0.052%
28	10.187	1199	1207	1215	rBV	1630246	2485939	36.91%	2.541%
29	10.257	1215	1219	1229	rVV9	11850	35276	0.52%	0.036%
30	10.340	1229	1233	1239	rVB4	15402	26648	0.40%	0.027%
31	10.552	1262	1269	1283	rBV	1976987	3439276	51.06%	3.515%
32	10.687	1285	1292	1314	rVB	599902	1105020	16.41%	1.129%
33	11.087	1355	1360	1372	rVB4	94337	167890	2.49%	0.172%
34	11.334	1397	1402	1411	rBV3	26943	66629	0.99%	0.068%
35	11.910	1494	1500	1505	rVV2	18483	35993	0.53%	0.037%
36	11.981	1508	1512	1518	rBV2	18460	26077	0.39%	0.027%

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Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP012925.MA.M
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37	12.046	1518	1523	1529	rBV2	28159	52150	0.77%	0.053%
38	12.140	1533	1539	1548	rVV3	26251	49474	0.73%	0.051%
39	12.763	1634	1645	1655	rVV	90834	172103	2.56%	0.176%
40	13.010	1681	1687	1695	rVB	107234	163427	2.43%	0.167%
41	13.228	1720	1724	1728	rVV5	29957	40422	0.60%	0.041%
42	13.310	1733	1738	1743	rVV	19398	32477	0.48%	0.033%
43	13.804	1815	1822	1827	rVV	2849087	3704945	55.01%	3.787%
44	13.846	1827	1829	1835	rVB	29408	43079	0.64%	0.044%
45	14.093	1860	1871	1883	rBV	3065007	4379676	65.03%	4.477%
46	14.398	1913	1923	1930	rBV	3045281	4693735	69.69%	4.798%
47	14.463	1930	1934	1941	rVB	69427	110034	1.63%	0.112%
48	14.622	1954	1961	1972	rBV	1062160	1758872	26.11%	1.798%
49	14.751	1979	1983	1988	rVB	83305	104495	1.55%	0.107%
50	14.804	1988	1992	2000	rVB2	29371	57102	0.85%	0.058%
51	15.228	2058	2064	2068	rVB2	22374	39410	0.59%	0.040%
52	15.404	2088	2094	2100	rBV	3529495	5311302	78.86%	5.429%
53	15.457	2100	2103	2108	rVB2	58167	78041	1.16%	0.080%
54	15.516	2108	2113	2122	rBV	850661	1202001	17.85%	1.229%
55	15.587	2122	2125	2129	rBV3	21445	28049	0.42%	0.029%
56	15.775	2151	2157	2166	rVB	2350177	3351769	49.76%	3.426%
57	16.151	2217	2221	2222	rBV4	22657	29200	0.43%	0.030%
58	16.175	2222	2225	2229	rVB	41995	54563	0.81%	0.056%
59	16.275	2238	2242	2245	rBV	40539	53857	0.80%	0.055%
60	16.357	2247	2256	2260	rVV2	313139	528679	7.85%	0.540%
61	16.392	2260	2262	2265	rVV2	58641	59767	0.89%	0.061%
62	16.434	2265	2269	2272	rVV2	79311	107101	1.59%	0.109%
63	16.481	2272	2277	2284	rVB4	67402	139792	2.08%	0.143%
64	16.592	2291	2296	2303	rBV6	85961	180190	2.68%	0.184%
65	16.663	2305	2308	2312	rVB2	29249	39475	0.59%	0.040%
66	16.839	2335	2338	2341	rBV	20588	28480	0.42%	0.029%
67	17.010	2364	2367	2370	rBV2	27873	43518	0.65%	0.044%
68	17.045	2370	2373	2379	rVB	70997	82691	1.23%	0.085%
69	17.192	2392	2398	2403	rBV2	3789869	5575539	82.78%	5.699%
70	17.239	2403	2406	2411	rVV	568456	819915	12.17%	0.838%
71	17.298	2411	2416	2428	rVV	3689985	5466904	81.17%	5.588%
72	17.386	2428	2431	2440	rVB7	36164	66847	0.99%	0.068%
73	17.586	2460	2465	2468	rBV2	143368	276142	4.10%	0.282%
74	17.904	2516	2519	2522	rBV5	31509	46896	0.70%	0.048%
75	18.004	2533	2536	2537	rBV2	27412	32609	0.48%	0.033%
76	18.098	2549	2552	2555	rBV	44703	57775	0.86%	0.059%

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 Title : SVOA CALIBRATION

77	18.139	2555	2559	2567	rBV3	614728	931080	13.82%	0.952%
78	18.210	2567	2571	2576	rVB	259574	363076	5.39%	0.371%
79	18.310	2584	2588	2592	rBV2	144345	242557	3.60%	0.248%
80	18.451	2608	2612	2617	rBV5	138915	221665	3.29%	0.227%
81	18.528	2622	2625	2629	rVB2	62384	74205	1.10%	0.076%
82	18.628	2638	2642	2645	rBV2	69333	104516	1.55%	0.107%
83	18.822	2673	2675	2680	rVB	74344	92811	1.38%	0.095%
84	19.016	2704	2708	2711	rBV2	117553	207391	3.08%	0.212%
85	19.051	2711	2714	2720	rVB4	232139	356781	5.30%	0.365%
86	19.110	2720	2724	2730	rBV2	157208	250276	3.72%	0.256%
87	19.328	2757	2761	2767	rBV	1128904	1571051	23.33%	1.606%
88	19.475	2782	2786	2792	rBV5	135747	250384	3.72%	0.256%
89	19.569	2798	2802	2809	rVB	548679	766460	11.38%	0.783%
90	19.675	2815	2820	2824	rBV	4548905	6214033	92.26%	6.351%
91	20.151	2898	2901	2905	rBV2	319462	369621	5.49%	0.378%
92	20.416	2943	2946	2952	rVB3	313092	384766	5.71%	0.393%
93	20.975	3038	3041	3044	rVB	364463	383670	5.70%	0.392%
94	21.327	3097	3101	3112	rVB2	1058696	1825650	27.11%	1.866%
95	21.580	3139	3144	3149	rBV	2641425	3741641	55.55%	3.824%
96	21.651	3149	3156	3161	rBV	4383902	6735315	100.00%	6.884%
97	22.586	3310	3315	3326	rVB2	533058	1570513	23.32%	1.605%
98	24.210	3587	3591	3598	rVB2	723319	1265073	18.78%	1.293%
99	24.804	3685	3692	3701	rBV	1321977	3333135	49.49%	3.407%
100	25.027	3722	3730	3741	rVB	2546063	6585934	97.78%	6.732%

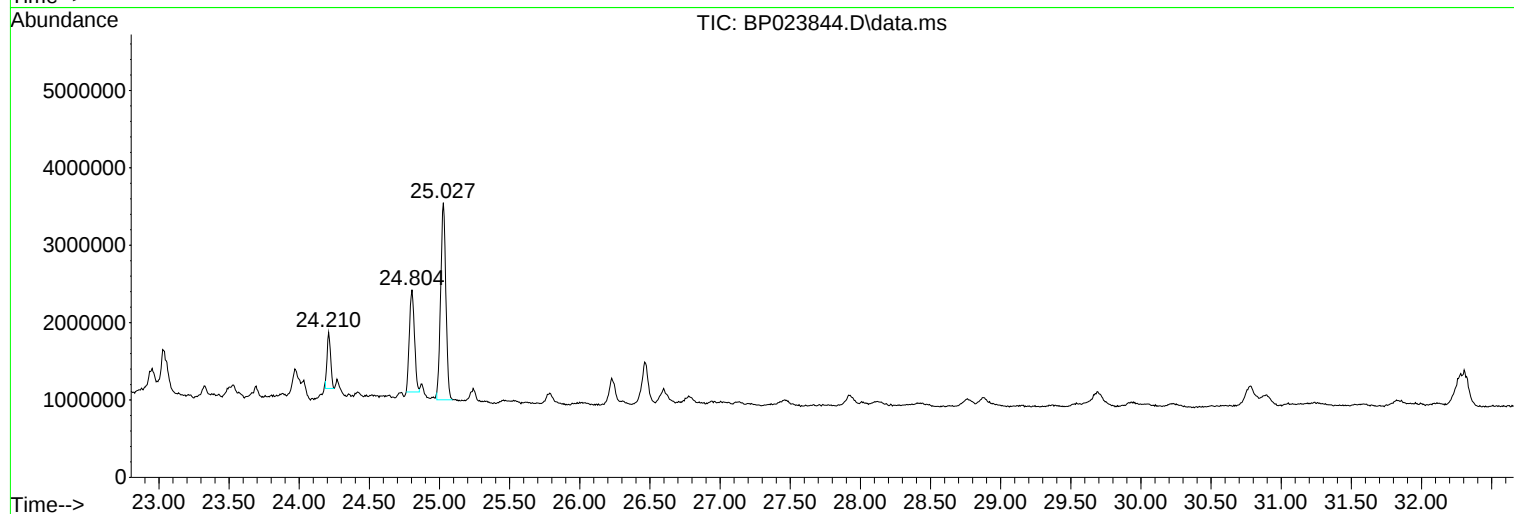
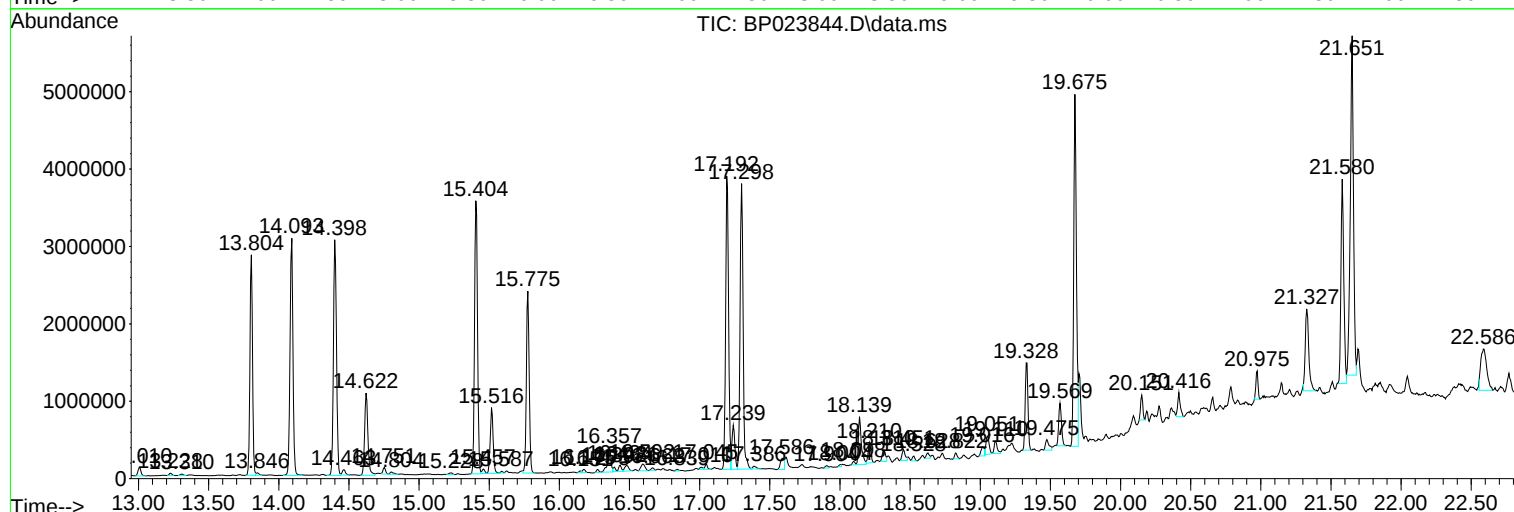
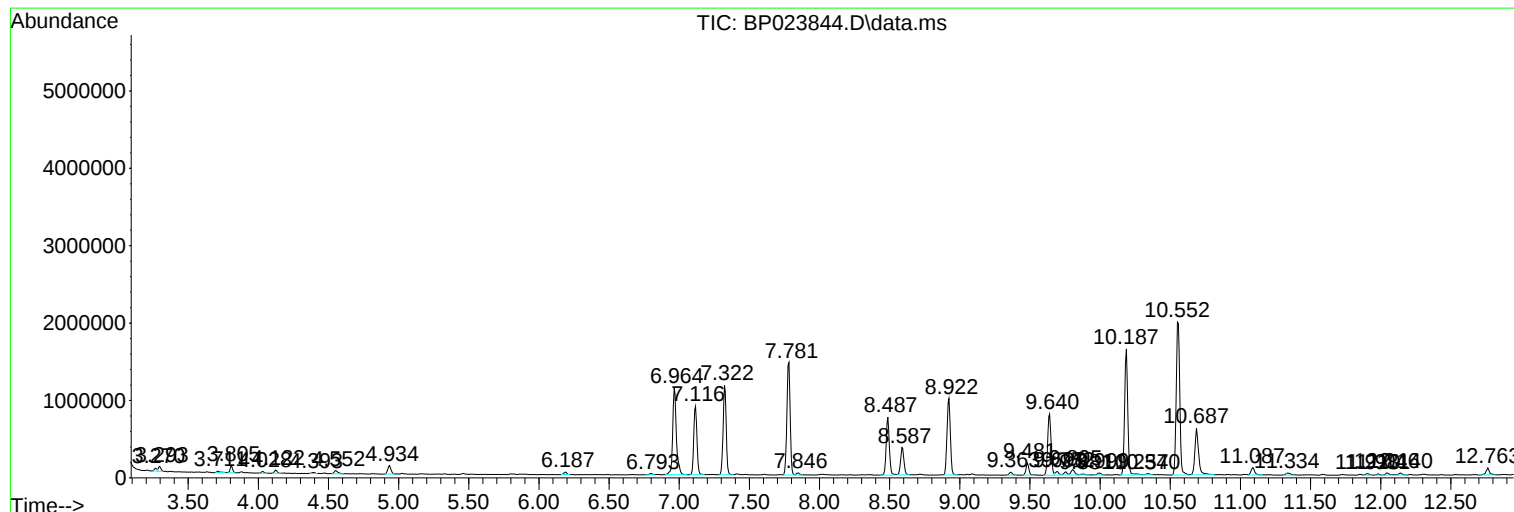
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ALS Vial : 28 Sample Multiplier: 1

Instrument :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP012925.MA.M
Quant Title : SVOA CALIBRATION

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



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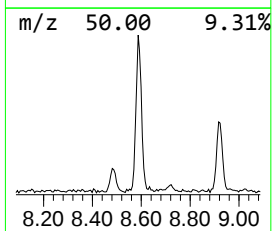
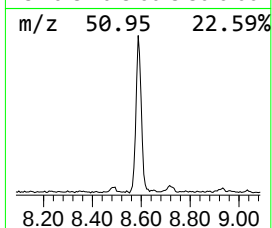
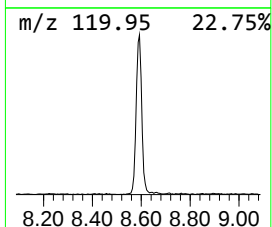
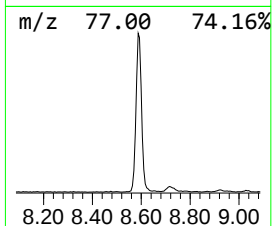
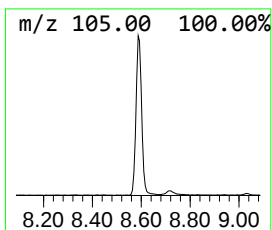
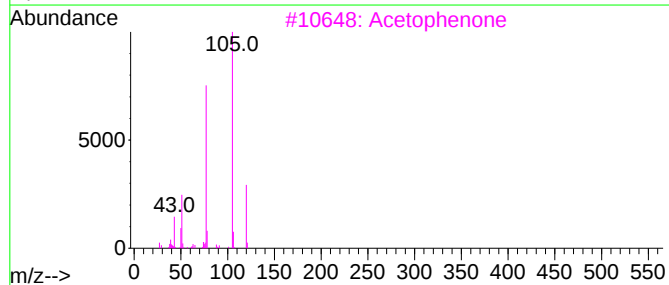
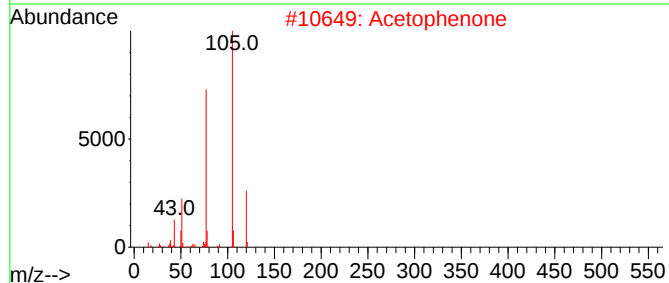
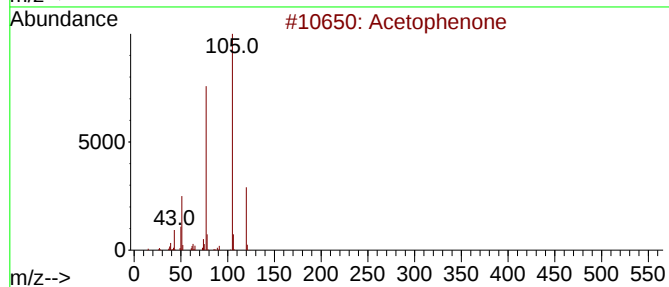
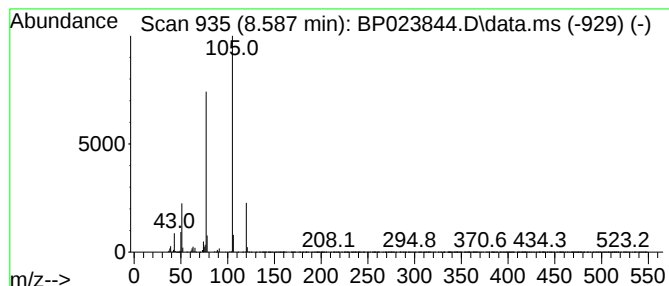
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP012925.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Aceto phenone Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.587	4.94 ng/ul	568145	1,4-Dichlorobenzene-d4	7.781

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Acetophenone			120	C8H8O	000098-86-2	95
2	Acetophenone			120	C8H8O	000098-86-2	94
3	Acetophenone			120	C8H8O	000098-86-2	94
4	Acetophenone			120	C8H8O	000098-86-2	94
5	Acetophenone			120	C8H8O	000098-86-2	91



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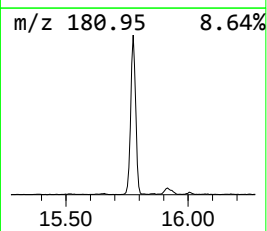
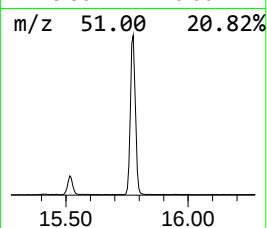
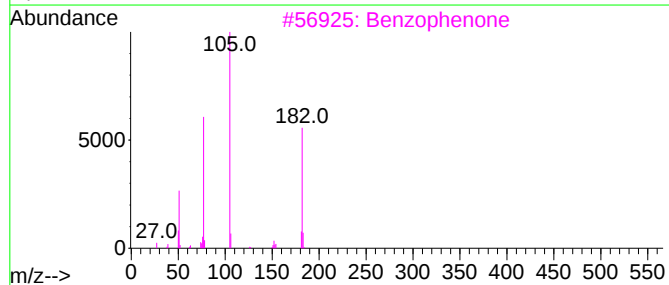
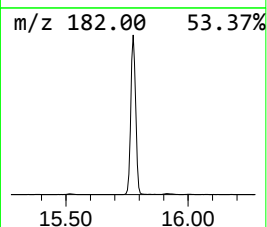
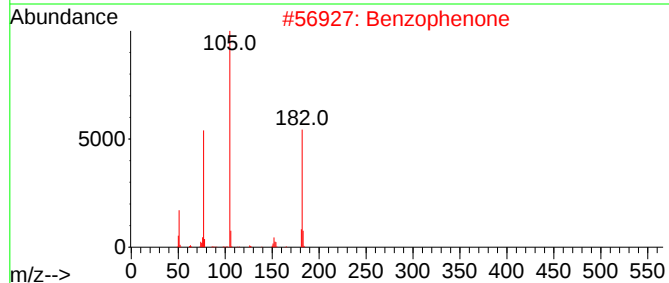
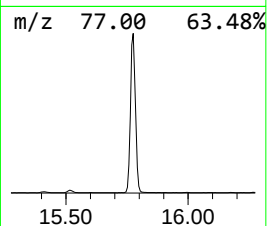
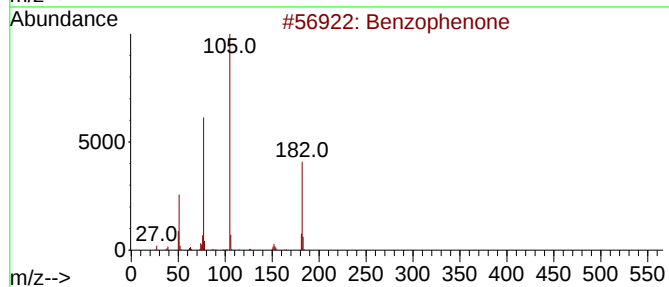
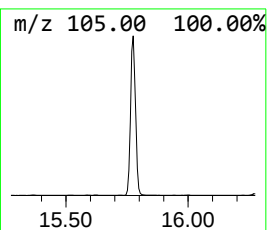
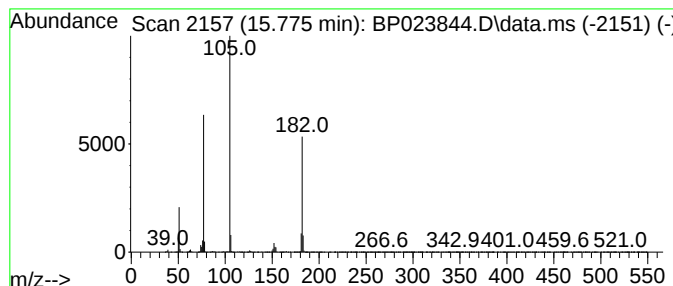
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP012925.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 Benzophenone Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.775	14.28 ng/ul	3351770	Acenaphthene-d10	14.398

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzophenone	182	C13H10O	000119-61-9	97
2			Benzophenone	182	C13H10O	000119-61-9	96
3			Benzophenone	182	C13H10O	000119-61-9	96
4			Benzophenone	182	C13H10O	000119-61-9	95
5			Benzophenone	182	C13H10O	000119-61-9	91



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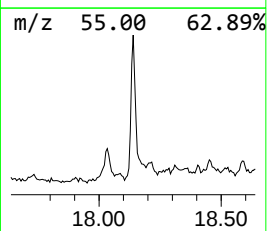
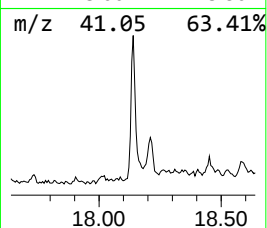
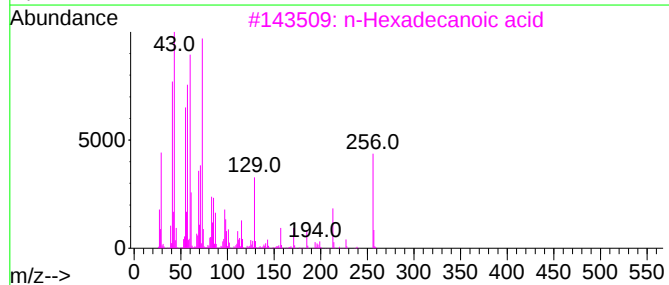
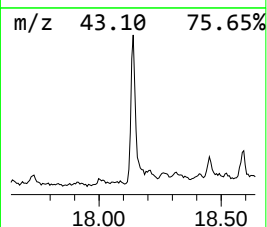
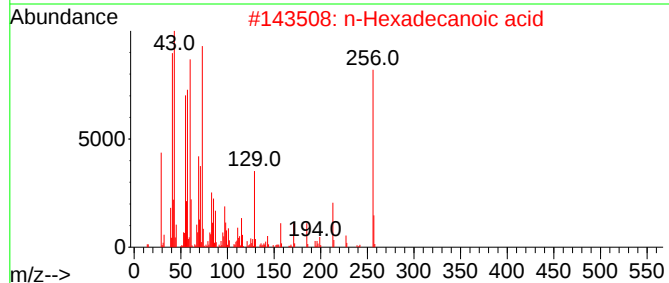
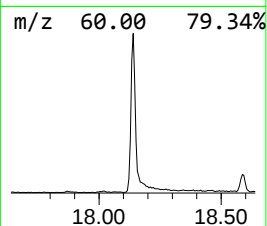
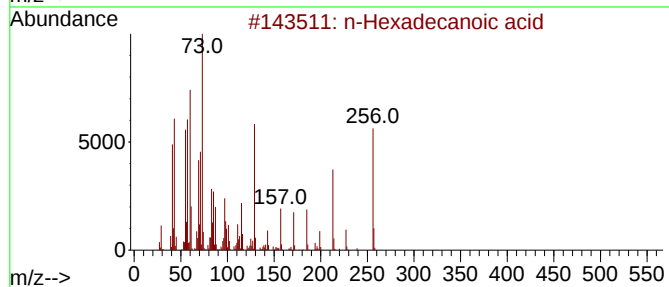
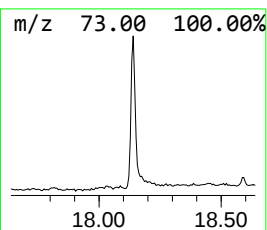
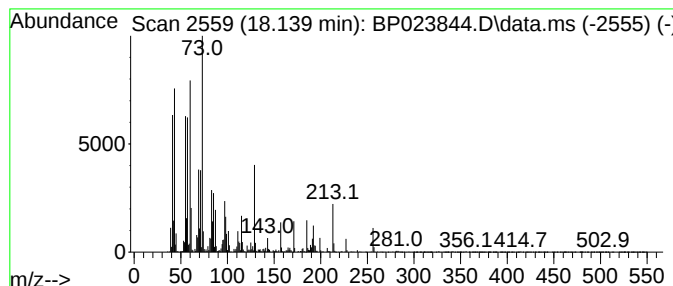
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP012925.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 n-Hexadecanoic acid Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.139	3.34 ng/ul	931080	Phenanthrene-d10	17.192

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	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP013125\
Data File : BP023844.D
Acq On : 01 Feb 2025 04:16
Operator : RC/JU
Sample : Q1224-08
Misc :
ALS Vial : 28 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A6306

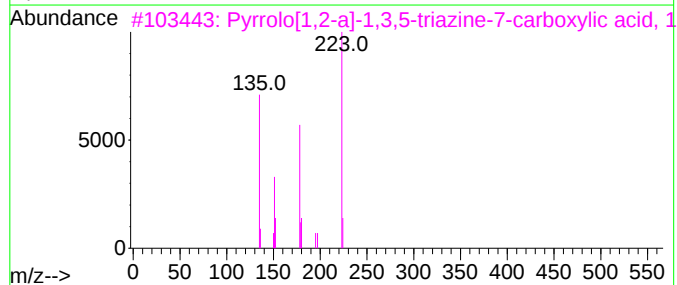
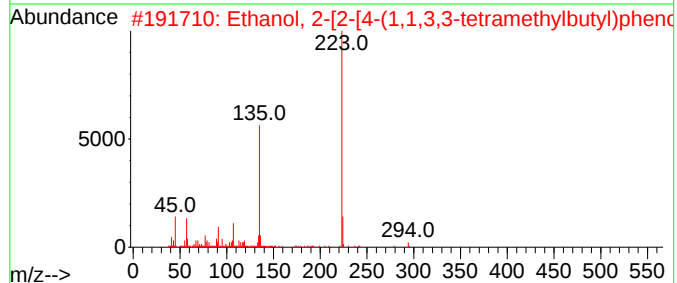
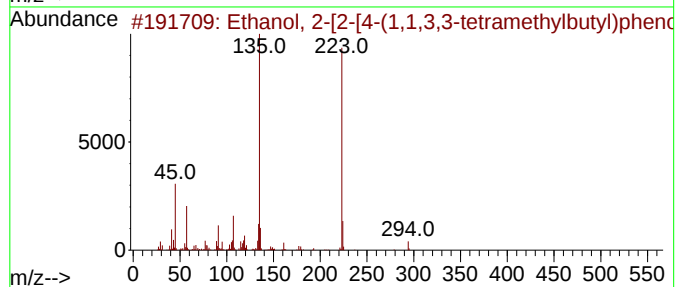
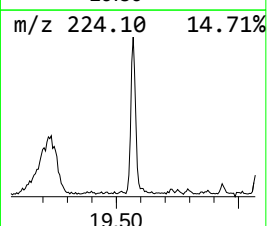
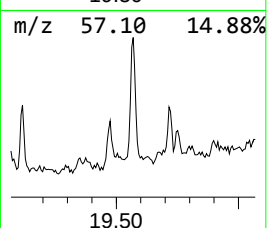
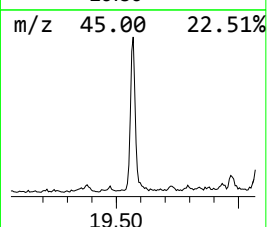
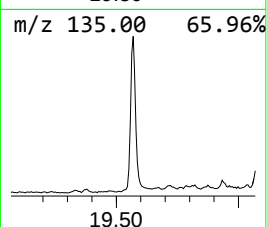
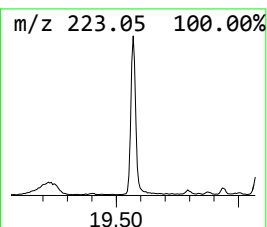
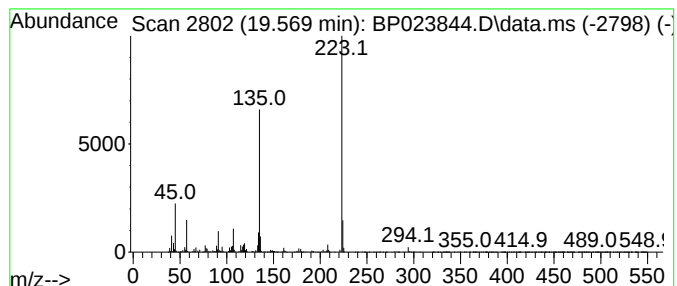
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 Ethanol, 2-[2-[4-(1,1,3,3-t... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.569	2.28 ng/ul	766460	Chrysene-d12	21.651

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Ethanol, 2-[2-[4-(1,1,3,3-tetram...	294	C18H30O3	002315-61-9	98
2			Ethanol, 2-[2-[4-(1,1,3,3-tetram...	294	C18H30O3	002315-61-9	93
3			Pyrrolo[1,2-a]-1,3,5-triazine-7-...	223	C9H9N3O4	054449-89-7	83
4			Glycine, N-(m-anisoyl)-, methyl ...	223	C11H13NO4	1000299-70-7	59
5			Ethanol, 2-[2-[4-(1,1,3,3-tetram...	294	C18H30O3	002315-61-9	43



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP013125\
Data File : BP023844.D
Acq On : 01 Feb 2025 04:16
Operator : RC/JU
Sample : Q1224-08
Misc :
ALS Vial : 28 Sample Multiplier: 1

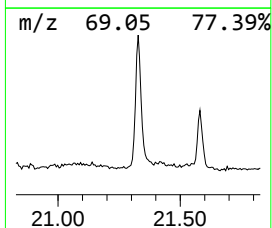
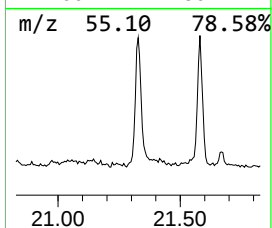
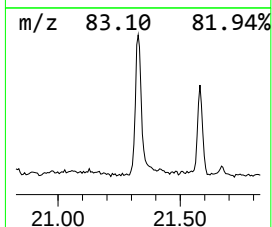
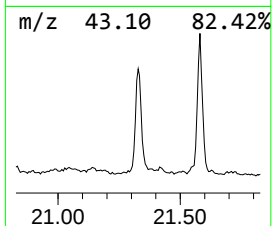
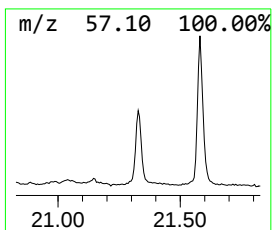
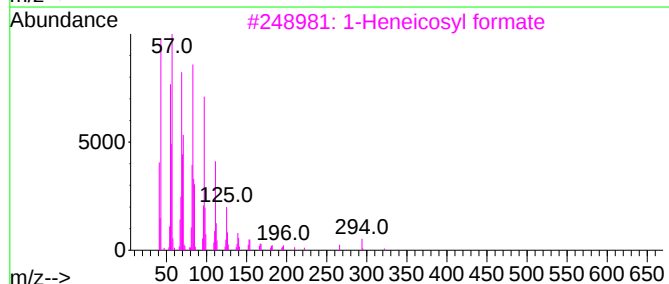
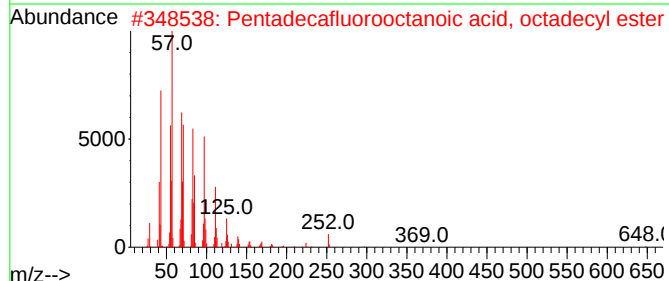
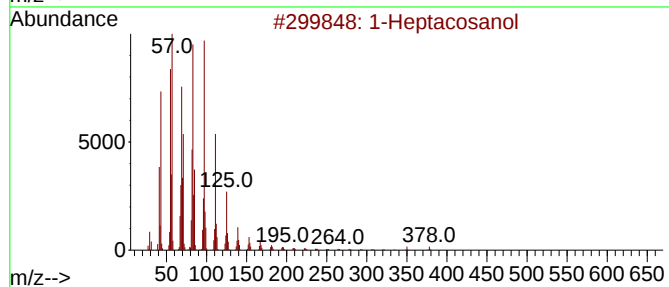
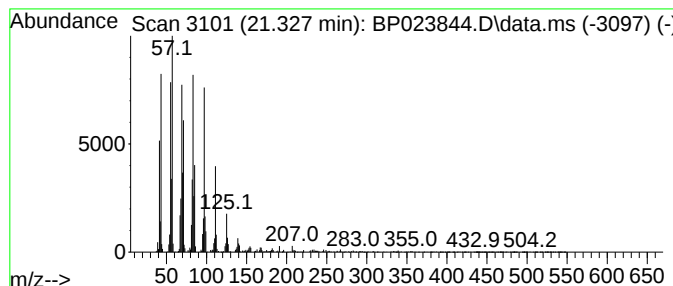
Instrument :
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ClientSampleId :
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 1-Heptacosanol Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.	
21.328	5.42 ng/ul	1825650	Chrysene-d12	21.651	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Heptacosanol	396	C27H56O	002004-39-9	94
2	Pentadecafluorooctanoic acid, octadecafluoro...	666	C26H37F15O2	1000406-04-8	94
3	1-Heneicosyl formate	340	C22H44O2	077899-03-7	92
4	Nonadecyl trifluoroacetate	380	C21H39F3O2	1000351-76-3	91
5	Trichloroacetic acid, hexadecyl ...	386	C18H33Cl3O2	074339-54-1	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP013125\
Data File : BP023844.D
Acq On : 01 Feb 2025 04:16
Operator : RC/JU
Sample : Q1224-08
Misc :
ALS Vial : 28 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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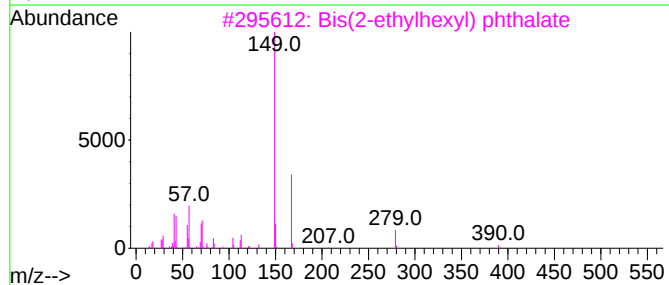
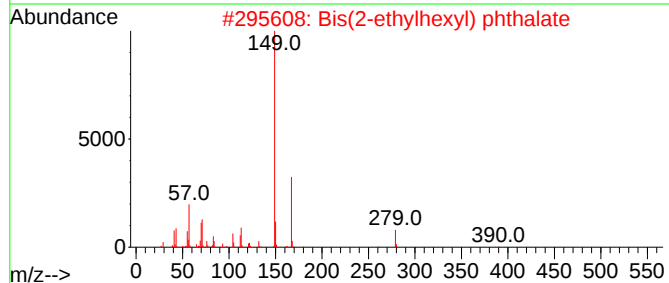
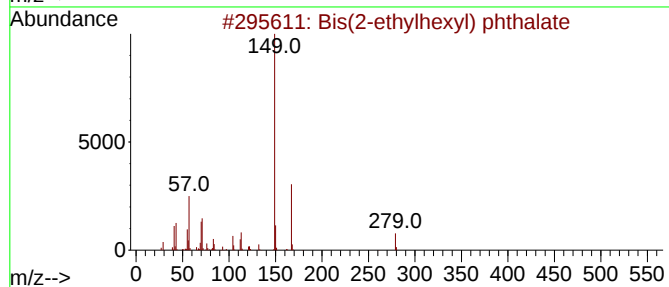
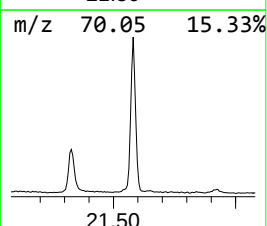
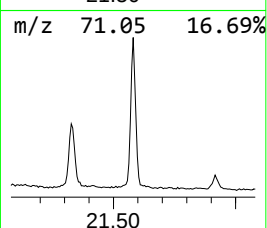
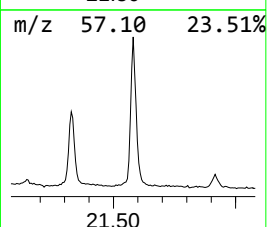
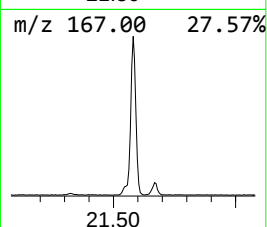
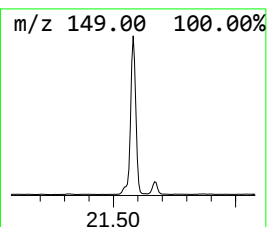
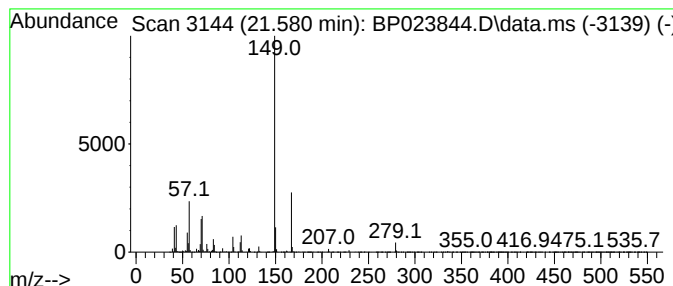
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 Bis(2-ethylhexyl) phthalate Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.580	11.11 ng/ul	3741640	Chrysene-d12	21.651

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Bis(2-ethylhexyl) phthalate	390	C24H38O4	000117-81-7	91
2			Bis(2-ethylhexyl) phthalate	390	C24H38O4	000117-81-7	91
3			Bis(2-ethylhexyl) phthalate	390	C24H38O4	000117-81-7	91
4			Phthalic acid, di(2-propylpentyl...)	390	C24H38O4	1000377-93-5	90
5			Phthalic acid, di(oct-3-yl) ester	390	C24H38O4	1000377-72-3	80



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP013125\
Data File : BP023844.D
Acq On : 01 Feb 2025 04:16
Operator : RC/JU
Sample : Q1224-08
Misc :
ALS Vial : 28 Sample Multiplier: 1

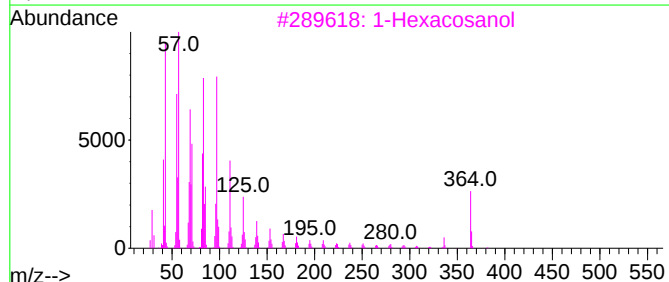
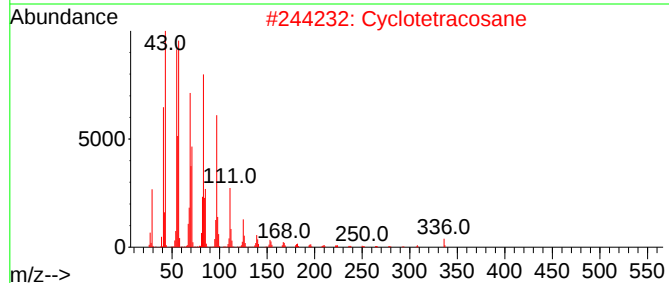
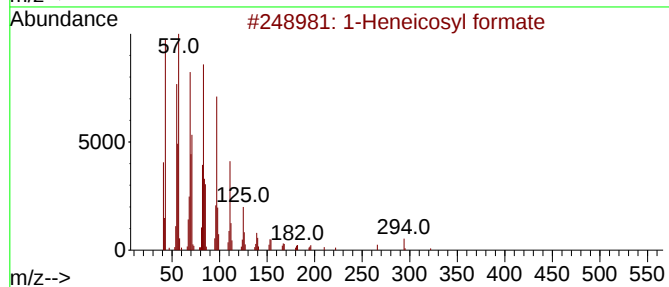
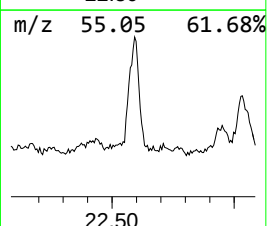
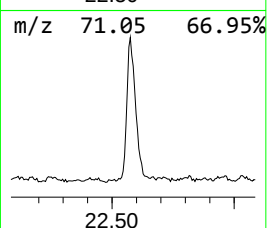
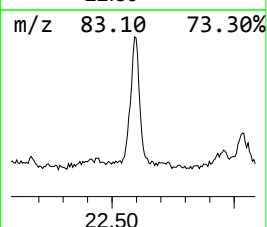
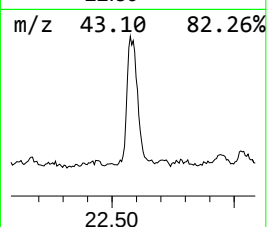
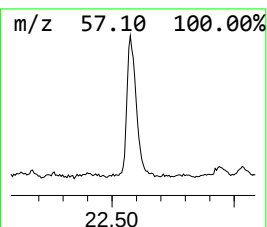
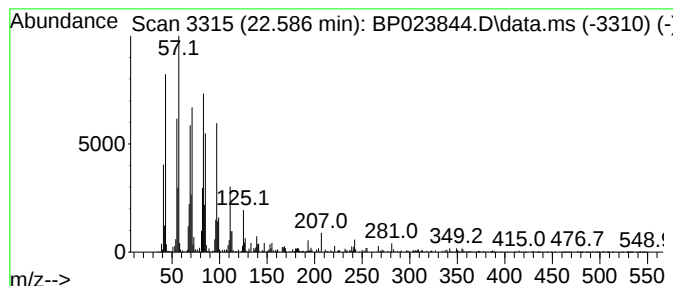
Instrument :
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ClientSampleId :
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 1-Heneicosyl formate Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.	
22.586	4.66 ng/ul	1570510	Chrysene-d12	21.651	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Heneicosyl formate	340	C22H44O2	077899-03-7	98
2	Cyclotetracosane	336	C24H48	000297-03-0	95
3	1-Hexacosanol	382	C26H54O	000506-52-5	93
4	1-Heneicosanol	312	C21H44O	015594-90-8	93
5	Eicosane, 9-cyclohexyl-	364	C26H52	004443-61-2	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP013125\
Data File : BP023844.D
Acq On : 01 Feb 2025 04:16
Operator : RC/JU
Sample : Q1224-08
Misc :
ALS Vial : 28 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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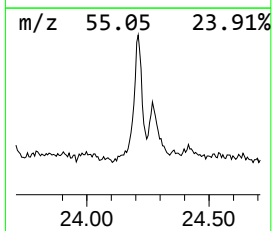
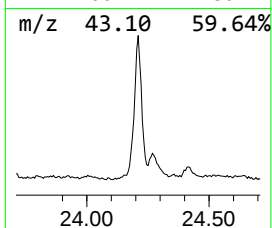
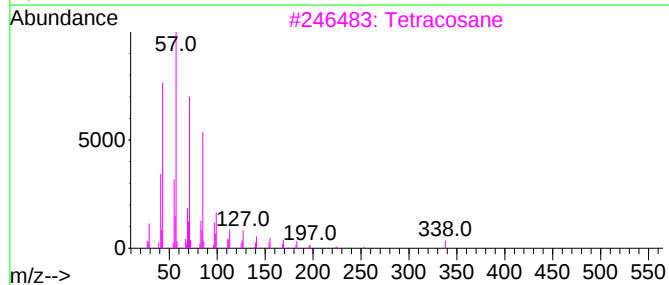
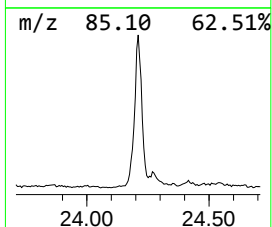
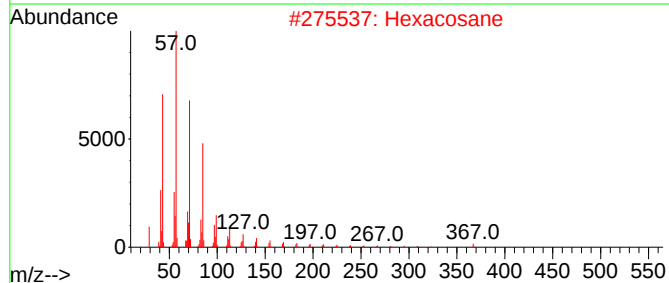
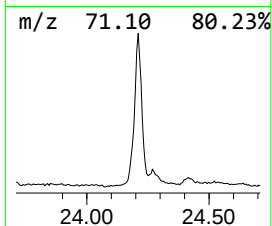
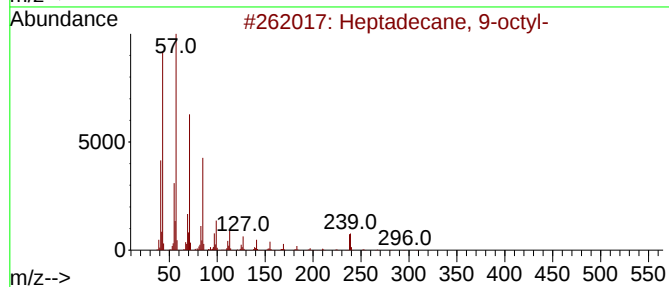
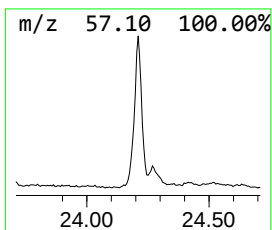
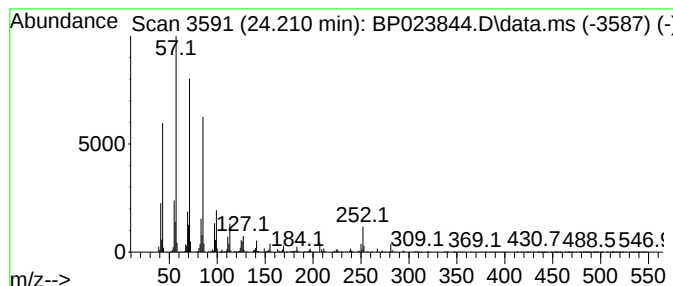
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 (DEL) Alkane: Straight-Chai... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.210	3.84 ng/ul	1265070	Perylene-d12	25.027

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Heptadecane, 9-octyl-	352	C25H52	007225-64-1	94
2		Hexacosane	366	C26H54	000630-01-3	93
3		Tetracosane	338	C24H50	000646-31-1	93
4		Eicosane, 9-octyl-	394	C28H58	013475-77-9	90
5		Heneicosane	296	C21H44	000629-94-7	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP013125\
Data File : BP023844.D
Acq On : 01 Feb 2025 04:16
Operator : RC/JU
Sample : Q1224-08
Misc :
ALS Vial : 28 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP012925.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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Benzophenone	15.775	14.3	ng/ul	3351770	3	14.398	4693740	20.0
n-Hexadecanoic ...	18.139	3.3	ng/ul	931080	4	17.192	5575540	20.0
Ethanol, 2-[2-[...	19.569	2.3	ng/ul	766460	5	21.651	6735320	20.0
1-Heptacosanol	21.328	5.4	ng/ul	1825650	5	21.651	6735320	20.0
Bis(2-ethylhexy...	21.580	11.1	ng/ul	3741640	5	21.651	6735320	20.0
1-Heneicosyl fo...	22.586	4.7	ng/ul	1570510	5	21.651	6735320	20.0
(DEL) Alkane: S...	24.210	3.8	ng/ul	1265070	6	25.027	6585930	20.0