

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP020325\
 Data File : BP023874.D
 Acq On : 03 Feb 2025 11:11
 Operator : RC/JU
 Sample : Q1248-01
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 A6318

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: BP023874.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.293	32	35	47	rVB	64523	98115	1.00%	0.085%
2	3.434	56	59	64	rVB3	21244	25342	0.26%	0.022%
3	3.717	104	107	117	rBV	84061	144180	1.47%	0.125%
4	4.028	156	160	164	rBV	27718	34259	0.35%	0.030%
5	4.675	266	270	276	rVB	33334	47987	0.49%	0.042%
6	6.164	518	523	530	rBV	61485	93310	0.95%	0.081%
7	6.793	623	630	634	rBV4	19799	35423	0.36%	0.031%
8	6.958	650	658	677	rVV	1168302	1950606	19.88%	1.697%
9	7.111	677	684	701	rVB	933452	1469019	14.97%	1.278%
10	7.316	713	719	729	rVV	1131648	1751055	17.85%	1.523%
11	7.775	785	797	807	rBV	1358003	2263220	23.07%	1.969%
12	7.916	817	821	828	rVB	14451	25698	0.26%	0.022%
13	8.016	828	838	844	rBV	30327	60952	0.62%	0.053%
14	8.481	910	917	927	rVV	1462874	2230708	22.74%	1.940%
15	8.922	985	992	1001	rVV	1039672	1660805	16.93%	1.445%
16	9.334	1056	1062	1065	rBV	42700	72429	0.74%	0.063%
17	9.481	1079	1087	1095	rVB3	33577	61927	0.63%	0.054%
18	9.640	1101	1114	1130	rBV	767323	1387625	14.14%	1.207%
19	9.875	1149	1154	1159	rVB8	14072	24184	0.25%	0.021%
20	9.975	1166	1171	1177	rVB8	12281	22531	0.23%	0.020%
21	10.181	1200	1206	1224	rVV	1400195	2182222	22.24%	1.898%
22	10.322	1224	1230	1237	rVB7	14214	35557	0.36%	0.031%
23	10.552	1259	1269	1274	rVV	1996005	3306816	33.71%	2.877%
24	10.604	1274	1278	1284	rVV	686287	1081203	11.02%	0.941%
25	10.681	1284	1291	1310	rVB	1200640	2023474	20.63%	1.760%
26	10.852	1315	1320	1326	rVB	41273	67049	0.68%	0.058%
27	11.087	1354	1360	1368	rVB7	18538	34326	0.35%	0.030%
28	11.604	1440	1448	1452	rBV3	13333	29045	0.30%	0.025%
29	11.981	1507	1512	1516	rVB2	22737	32388	0.33%	0.028%
30	12.134	1529	1538	1545	rVB	21435	42100	0.43%	0.037%
31	12.263	1556	1560	1564	rBV	18949	29661	0.30%	0.026%
32	13.134	1699	1708	1710	rBV9	14138	27539	0.28%	0.024%
33	13.222	1718	1723	1727	rBV	44602	62954	0.64%	0.055%
34	13.304	1732	1737	1743	rVV	49873	91061	0.93%	0.079%
35	13.369	1743	1748	1759	rVV6	17489	45200	0.46%	0.039%
36	13.675	1795	1800	1804	rBV6	39831	69690	0.71%	0.061%

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Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP012925.MA.M
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37	13.798	1814	1821	1830	rBV	2566206	3189285	32.51%	2.774%
38	14.081	1860	1869	1873	rBV2	2662413	4384288	44.69%	3.814%
39	14.298	1903	1906	1911	rVB2	19644	28131	0.29%	0.024%
40	14.393	1913	1922	1928	rBV	3096787	4439489	45.25%	3.862%
41	14.457	1928	1933	1940	rVV	1962470	2917136	29.73%	2.538%
42	14.522	1940	1944	1951	rVB	304882	389819	3.97%	0.339%
43	14.604	1951	1958	1973	rBV	1327929	1866065	19.02%	1.623%
44	14.722	1973	1978	1987	rVV6	57938	106137	1.08%	0.092%
45	14.822	1988	1995	1999	rVB3	46494	74967	0.76%	0.065%
46	14.951	2013	2017	2021	rVB4	16271	24579	0.25%	0.021%
47	15.198	2054	2059	2066	rBV7	22158	53128	0.54%	0.046%
48	15.263	2066	2070	2076	rVB2	21388	28105	0.29%	0.024%
49	15.398	2086	2093	2098	rBV	4013762	5460926	55.66%	4.750%
50	15.451	2098	2102	2107	rVV	1195086	1622280	16.54%	1.411%
51	15.510	2107	2112	2121	rVB	830939	1166509	11.89%	1.015%
52	15.651	2131	2136	2142	rVB7	19243	34679	0.35%	0.030%
53	15.769	2150	2156	2164	rVB	1261984	1646062	16.78%	1.432%
54	15.898	2175	2178	2182	rVB	18024	25230	0.26%	0.022%
55	16.145	2215	2220	2222	rBV4	31135	43795	0.45%	0.038%
56	16.345	2248	2254	2264	rBV2	461641	715625	7.29%	0.623%
57	16.587	2291	2295	2303	rVB4	33737	56900	0.58%	0.049%
58	16.728	2316	2319	2324	rVB2	25199	33446	0.34%	0.029%
59	16.851	2333	2340	2346	rVB2	156446	224766	2.29%	0.196%
60	16.922	2347	2352	2356	rBV	77680	123802	1.26%	0.108%
61	16.992	2356	2364	2374	rVV3	179433	451835	4.61%	0.393%
62	17.092	2378	2381	2385	rBV6	19543	33250	0.34%	0.029%
63	17.139	2386	2389	2391	rVV	42702	49196	0.50%	0.043%
64	17.186	2391	2397	2401	rVV	3724821	5310584	54.13%	4.620%
65	17.234	2401	2405	2410	rVV	2372579	3620358	36.90%	3.149%
66	17.292	2410	2415	2418	rVV	4906275	6859633	69.92%	5.967%
67	17.328	2418	2421	2428	rVB	2445804	3312796	33.77%	2.882%
68	17.504	2447	2451	2458	rVB2	36272	60925	0.62%	0.053%
69	17.628	2465	2472	2477	rBV3	95722	150717	1.54%	0.131%
70	17.675	2477	2480	2484	rVB4	19922	22863	0.23%	0.020%
71	17.722	2485	2488	2491	rVV4	33280	41623	0.42%	0.036%
72	17.763	2491	2495	2498	rVV3	60025	87189	0.89%	0.076%
73	17.828	2502	2506	2511	rVV3	107707	167671	1.71%	0.146%
74	17.898	2513	2518	2527	rVB	632861	878971	8.96%	0.765%
75	18.128	2552	2557	2563	rBV	444078	638435	6.51%	0.555%
76	18.351	2593	2595	2599	rVB	58852	64704	0.66%	0.056%

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77	18.433	2603	2609	2617	rVV5	144564	354275	3.61%	0.308%
78	18.510	2617	2622	2632	rVV2	191900	459306	4.68%	0.400%
79	18.598	2632	2637	2641	rVV4	112111	218021	2.22%	0.190%
80	18.651	2641	2646	2652	rVB	312352	441871	4.50%	0.384%
81	18.875	2680	2684	2687	rBV	124383	149542	1.52%	0.130%
82	18.910	2688	2690	2694	rVB3	42085	45658	0.47%	0.040%
83	18.998	2700	2705	2707	rBV2	199488	343011	3.50%	0.298%
84	19.028	2707	2710	2716	rVB	229822	355027	3.62%	0.309%
85	19.110	2721	2724	2728	rVV2	317122	365329	3.72%	0.318%
86	19.210	2738	2741	2744	rBV2	84453	107597	1.10%	0.094%
87	19.257	2745	2749	2752	rVV	292432	333402	3.40%	0.290%
88	19.310	2754	2758	2764	rVB	2117509	2574759	26.24%	2.240%
89	19.457	2780	2783	2786	rBV2	93201	134174	1.37%	0.117%
90	19.663	2812	2818	2827	rBV2	6673593	9810744	100.00%	8.534%
91	20.398	2939	2943	2948	rVB	844202	1042739	10.63%	0.907%
92	20.698	2991	2994	2997	rVB	338104	357778	3.65%	0.311%
93	21.322	3096	3100	3109	rVB	849795	1258096	12.82%	1.094%
94	21.480	3123	3127	3133	rVB	796839	1069160	10.90%	0.930%
95	21.639	3147	3154	3158	rBV	3789024	5998658	61.14%	5.218%
96	21.680	3158	3161	3167	rVB	696623	1019893	10.40%	0.887%
97	22.769	3343	3346	3353	rVB	567969	924981	9.43%	0.805%
98	23.933	3540	3544	3551	rBV	382724	764586	7.79%	0.665%
99	24.780	3679	3688	3695	rBV	3250751	7831679	79.83%	6.813%
100	25.010	3719	3727	3746	rVB	2157703	5972204	60.87%	5.195%

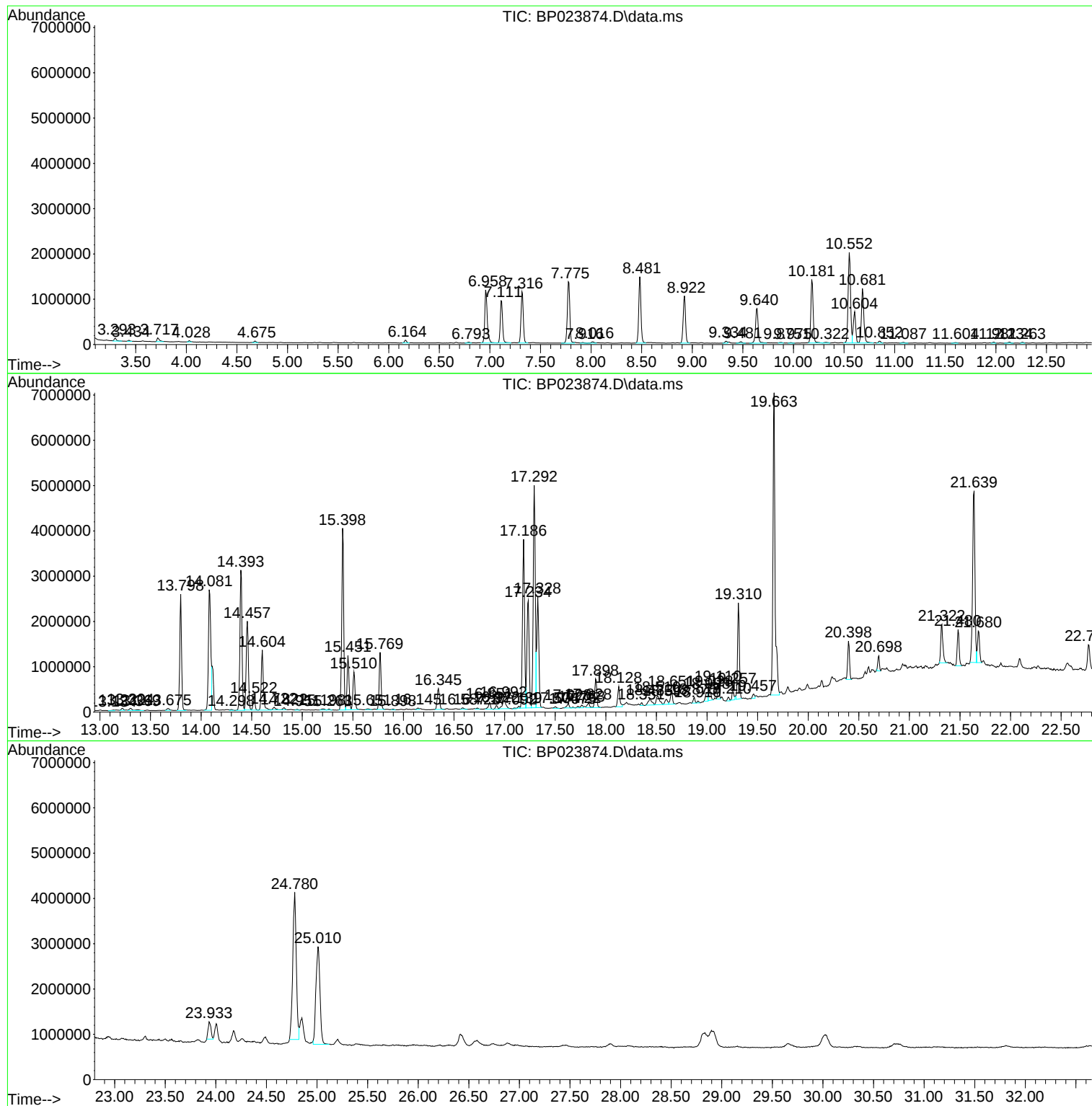
Sum of corrected areas: 114958049

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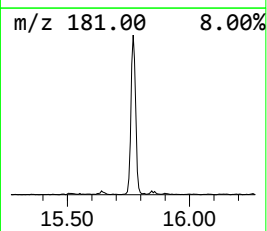
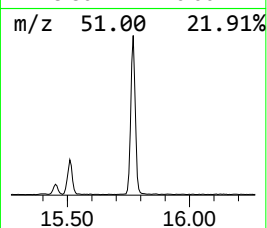
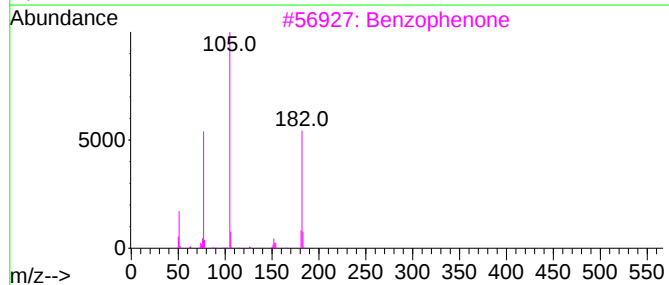
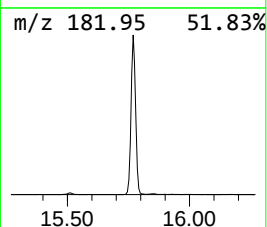
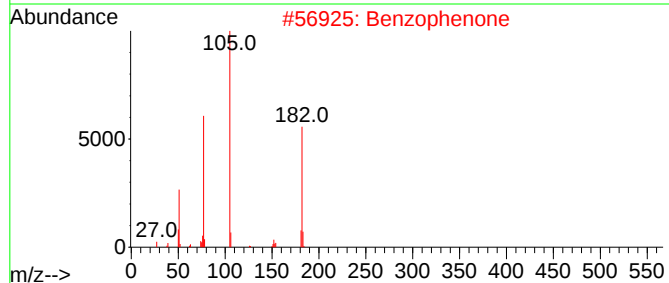
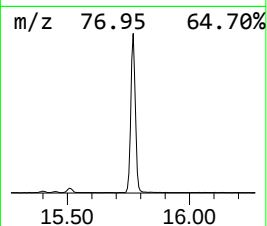
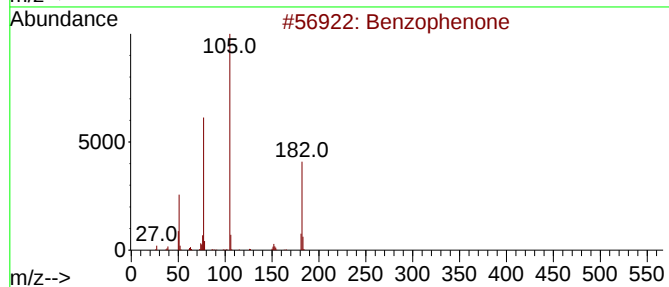
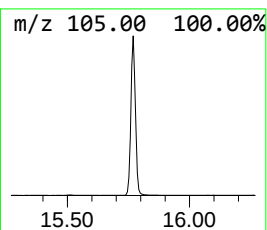
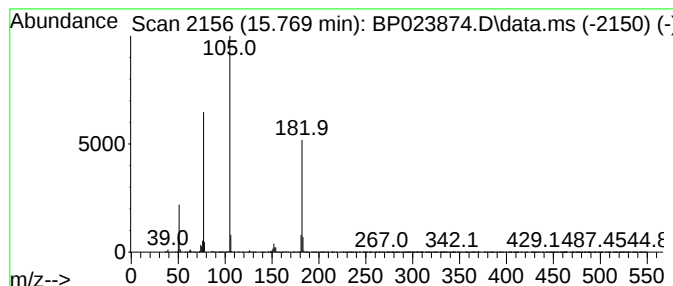
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP012925.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Benzophenone Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.769	7.42 ng/ul	1646060	Acenaphthene-d10	14.393

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	93
5			Benzophenone	182	C13H10O	000119-61-9	91



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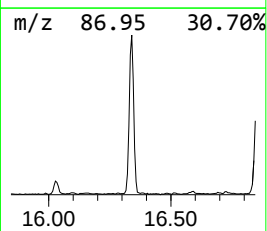
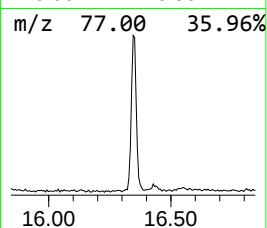
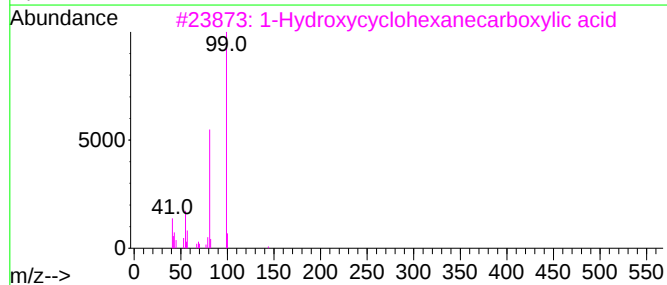
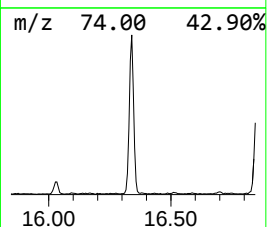
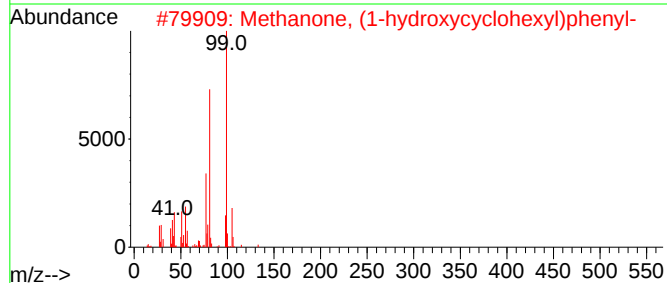
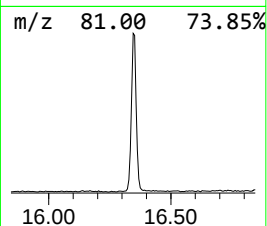
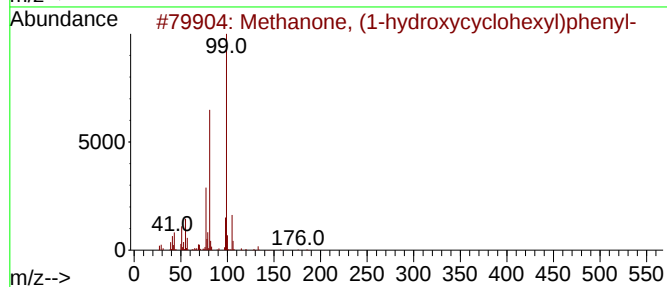
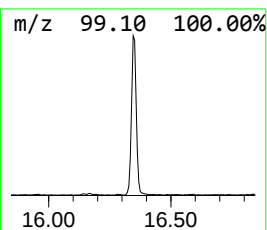
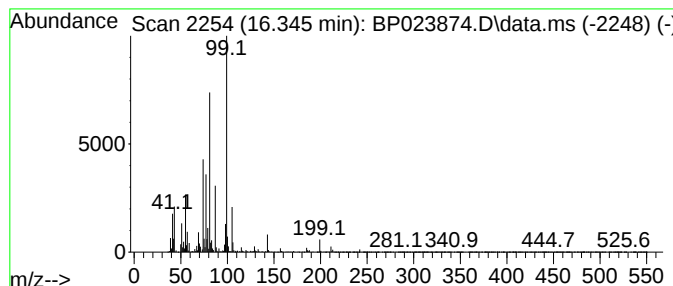
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Quant Title : SVOA CALIBRATION

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

Peak Number 2 Methanone, (1-hydroxycyclohexyl) Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.345	2.70 ng/ul	715625	Phenanthrene-d10	17.186

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Methanone, (1-hydroxycyclohexyl)...	204	C13H16O2	000947-19-3	93
2			Methanone, (1-hydroxycyclohexyl)...	204	C13H16O2	000947-19-3	58
3			1-Hydroxycyclohexanecarboxylic acid	144	C7H12O3	001123-28-0	46
4			Cyclohexanol, 1-ethyl-	128	C8H16O	001940-18-7	43
5			Cyclohexanol, 1-(aminomethyl)-	129	C7H15NO	004000-72-0	38



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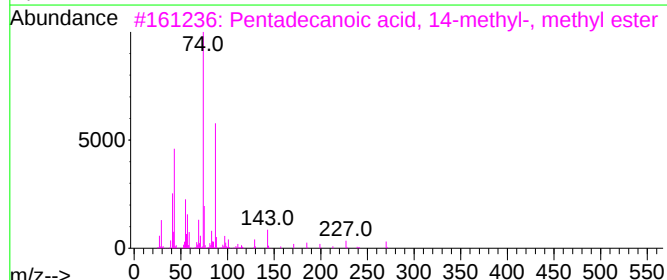
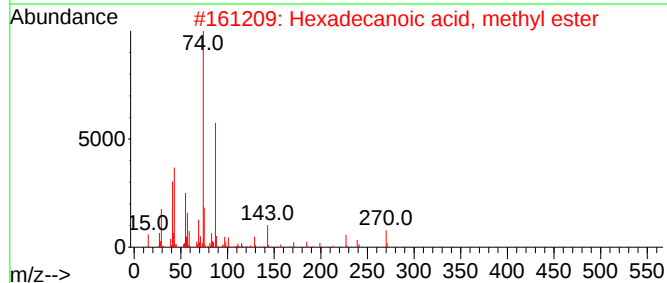
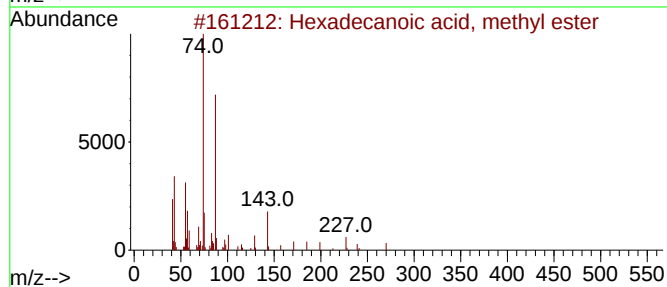
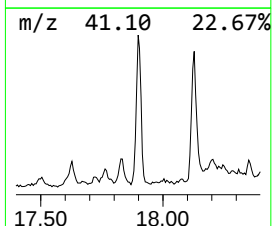
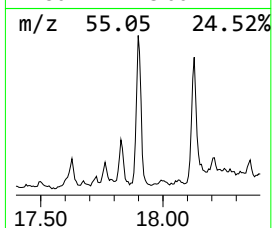
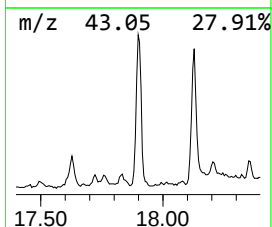
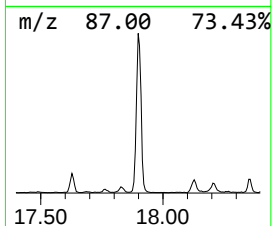
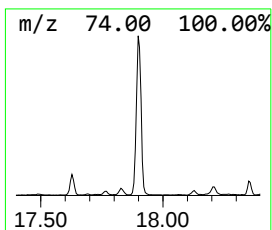
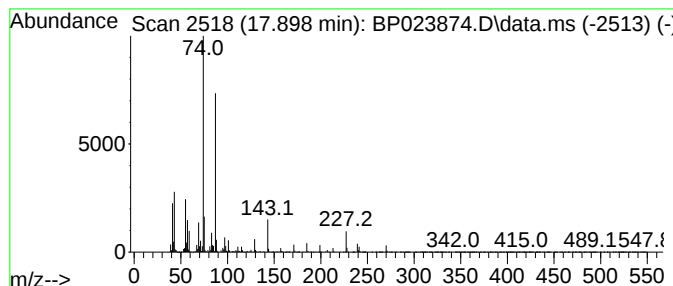
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Quant Title : SVOA CALIBRATION

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

Peak Number 3 Hexadecanoic acid, methyl e... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.898	3.31 ng/ul	878971	Phenanthrene-d10	17.186

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexadecanoic acid, methyl ester	270	C17H34O2	000112-39-0	98
2			Hexadecanoic acid, methyl ester	270	C17H34O2	000112-39-0	97
3			Pentadecanoic acid, 14-methyl-, ...	270	C17H34O2	005129-60-2	96
4			Tridecanoic acid, methyl ester	228	C14H28O2	001731-88-0	93
5			Methyl tetradecanoate	242	C15H30O2	000124-10-7	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP020325\
Data File : BP023874.D
Acq On : 03 Feb 2025 11:11
Operator : RC/JU
Sample : Q1248-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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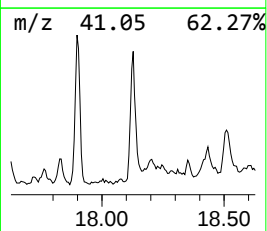
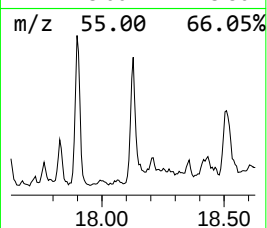
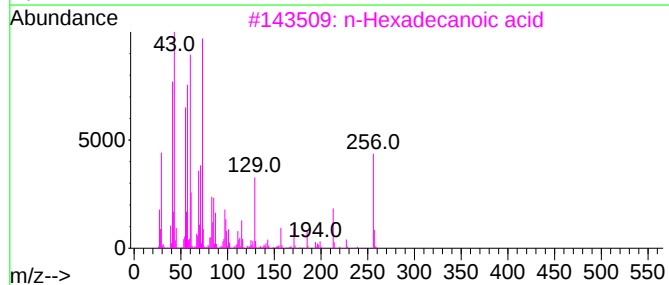
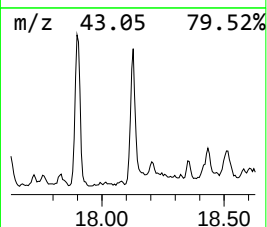
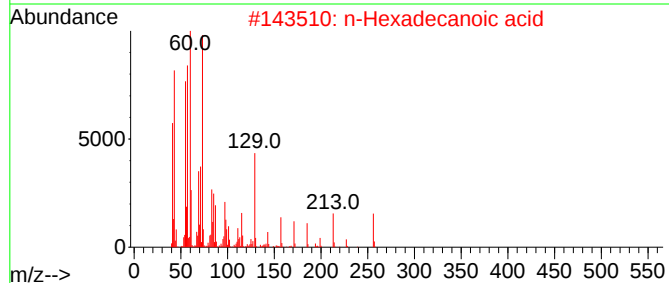
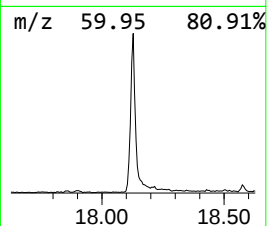
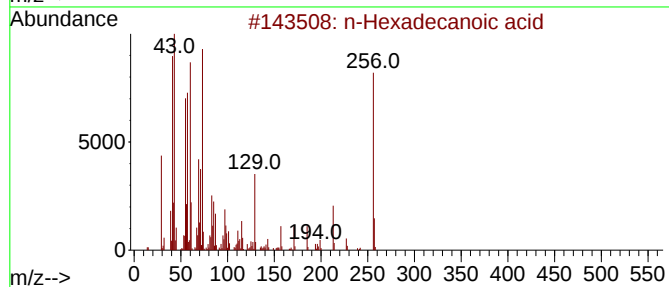
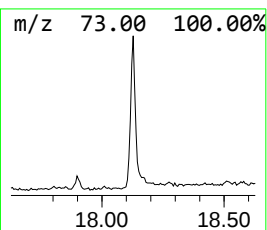
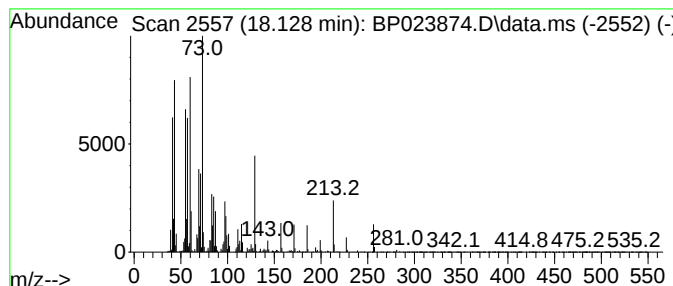
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 n-Hexadecanoic acid Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.128	2.40 ng/ul	638435	Phenanthrene-d10	17.186

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			Tridecanoic acid	214	C13H26O2	000638-53-9	89



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP020325\
Data File : BP023874.D
Acq On : 03 Feb 2025 11:11
Operator : RC/JU
Sample : Q1248-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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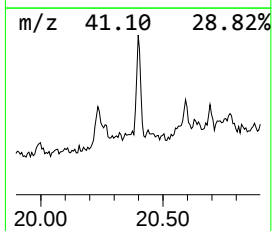
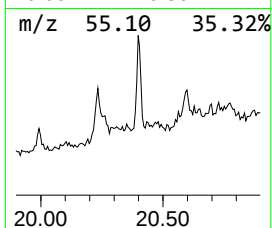
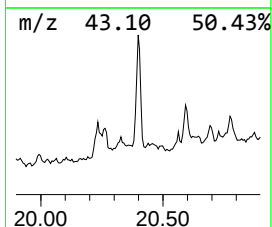
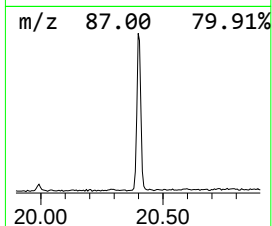
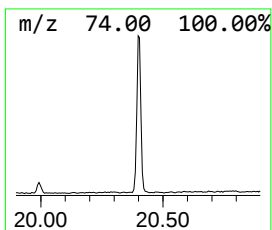
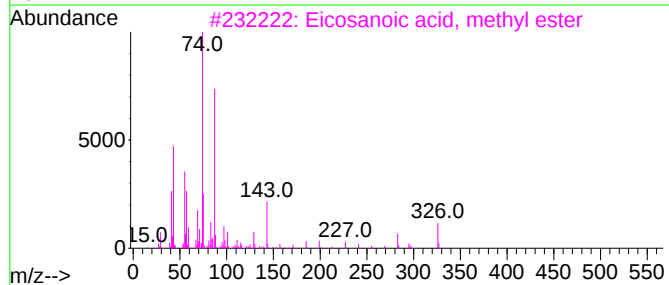
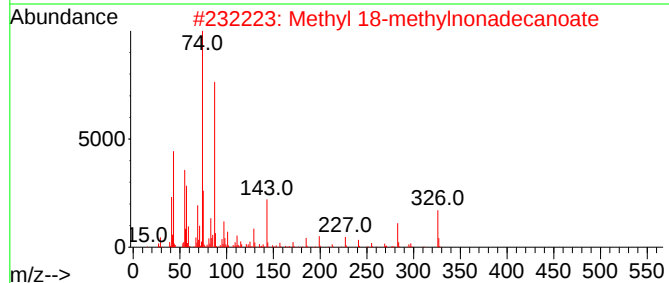
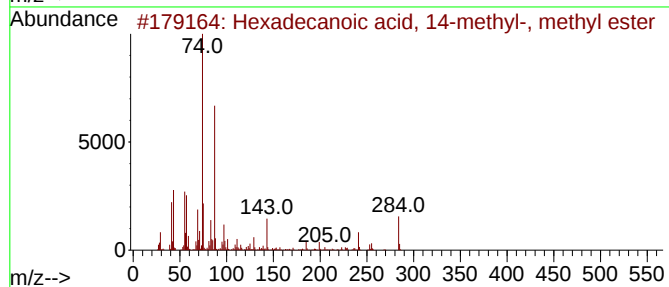
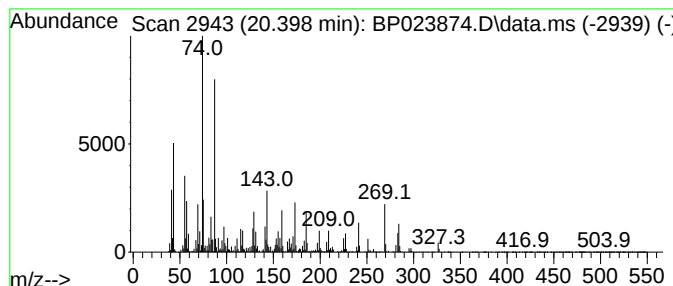
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 Hexadecanoic acid, 14-methy... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.398	3.48 ng/ul	1042740	Chrysene-d12	21.639

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexadecanoic acid, 14-methyl-, m...	284	C18H36O2	002490-49-5	98
2			Methyl 18-methylnonadecanoate	326	C21H42O2	1000424-52-4	98
3			Eicosanoic acid, methyl ester	326	C21H42O2	001120-28-1	96
4			1-Phenanthrenecarboxaldehyde, 1,...	284	C20H28O	013601-88-2	92
5			Hexadecanoic acid, 15-methyl-, m...	284	C18H36O2	006929-04-0	89



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP020325\
Data File : BP023874.D
Acq On : 03 Feb 2025 11:11
Operator : RC/JU
Sample : Q1248-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

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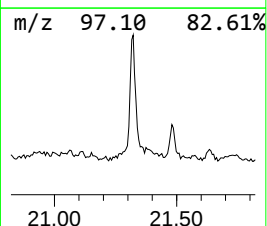
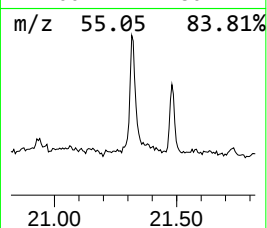
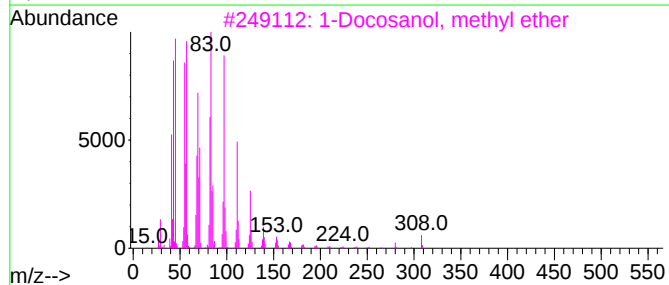
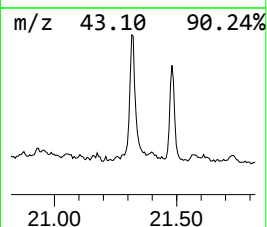
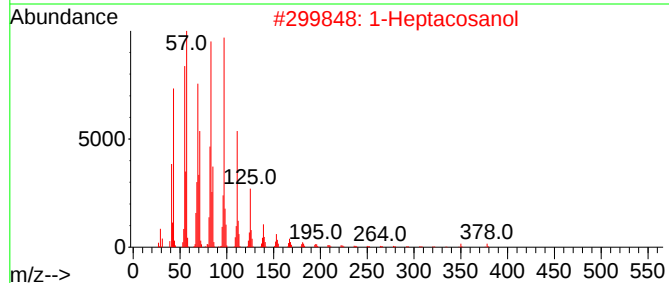
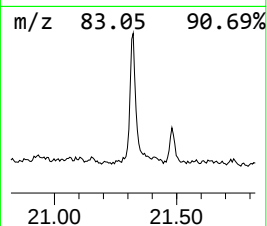
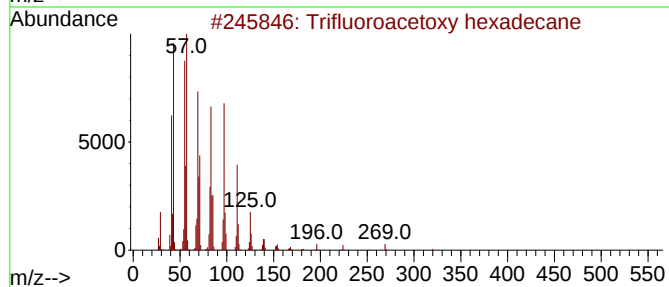
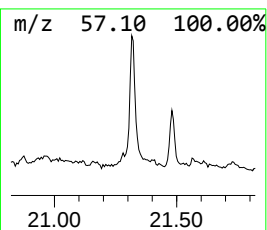
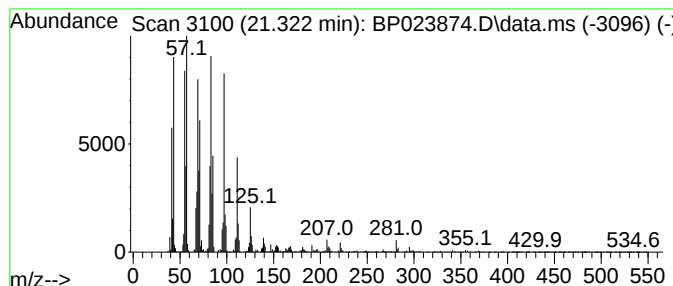
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 Trifluoroacetoxy hexadecane Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.322	4.19 ng/ul	1258100	Chrysene-d12	21.639

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Trifluoroacetoxy hexadecane	338	C18H33F3O2	006222-03-3	94
2			1-Heptacosanol	396	C27H56O	002004-39-9	94
3			1-Docosanol, methyl ether	340	C23H48O	1000333-91-9	93
4			1-Tetracosene	336	C24H48	010192-32-2	93
5			1-Heneicosanol	312	C21H44O	015594-90-8	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP020325\
Data File : BP023874.D
Acq On : 03 Feb 2025 11:11
Operator : RC/JU
Sample : Q1248-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

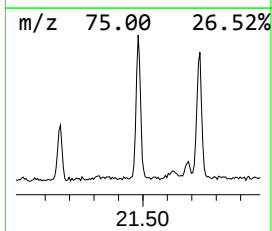
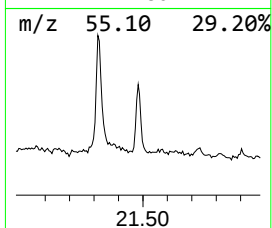
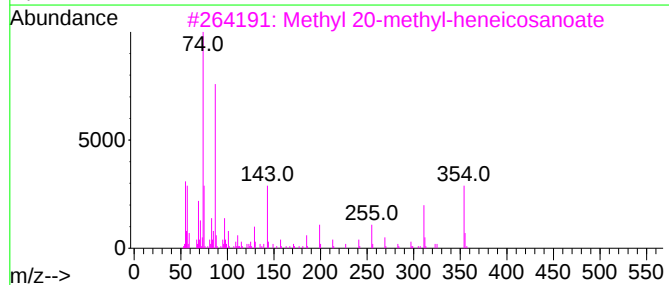
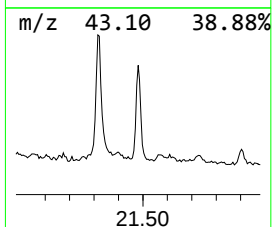
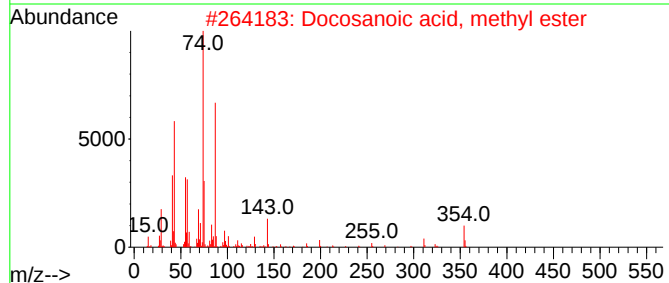
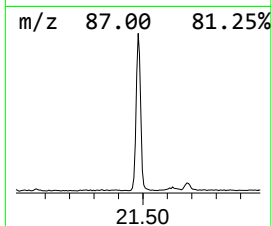
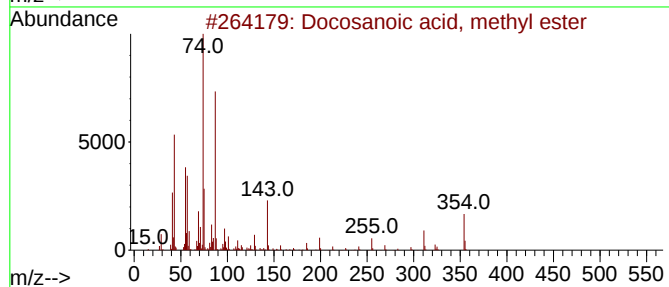
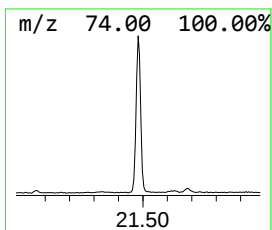
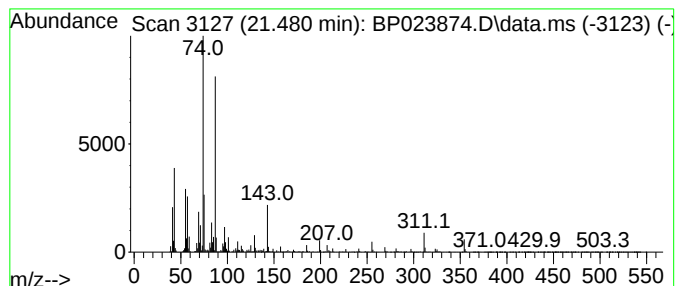
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 Docosanoic acid, methyl ester Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.	
21.480	3.56 ng/ul	1069160	Chrysene-d12	21.639	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Docosanoic acid, methyl ester	354	C23H46O2	000929-77-1	99
2	Docosanoic acid, methyl ester	354	C23H46O2	000929-77-1	98
3	Methyl 20-methyl-heneicosanoate	354	C23H46O2	1000336-47-4	98
4	Docosanoic acid, methyl ester	354	C23H46O2	000929-77-1	95
5	Docosanoic acid, methyl ester	354	C23H46O2	000929-77-1	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP020325\
Data File : BP023874.D
Acq On : 03 Feb 2025 11:11
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Sample : Q1248-01
Misc :
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Instrument :
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ClientSampleId :
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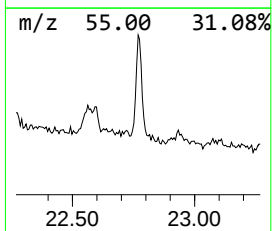
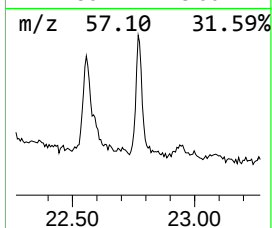
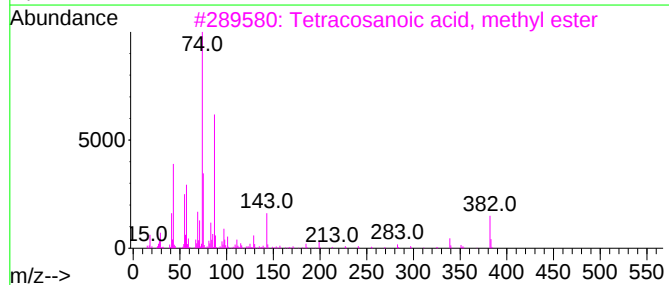
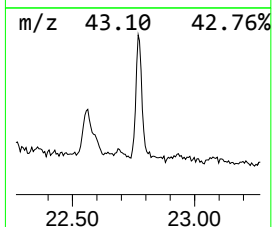
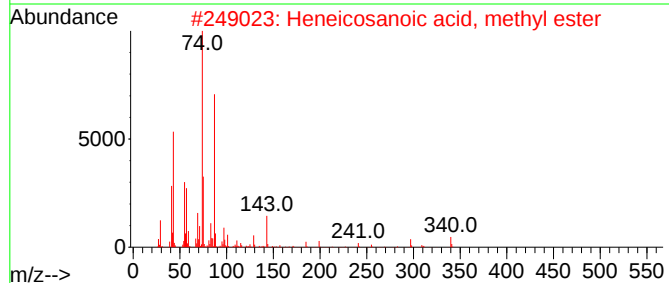
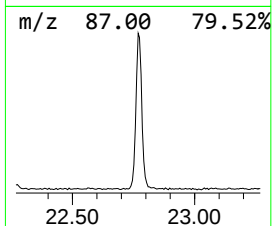
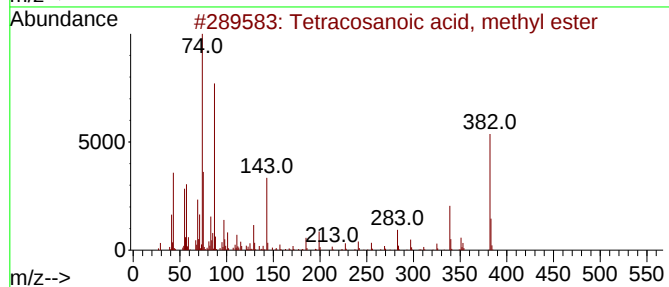
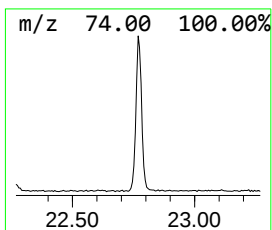
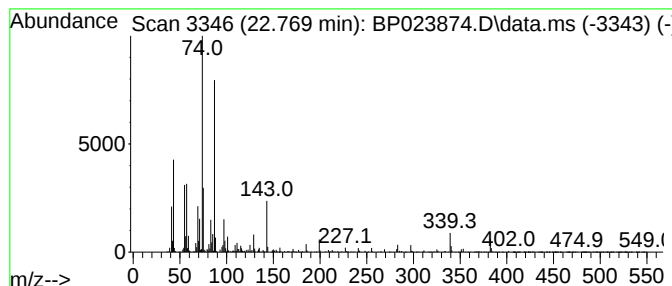
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 Tetracosanoic acid, methyl ... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.769	3.08 ng/ul	924981	Chrysene-d12	21.639

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tetracosanoic acid, methyl ester	382	C25H50O2	002442-49-1	98
2			Heneicosanoic acid, methyl ester	340	C22H44O2	006064-90-0	97
3			Tetracosanoic acid, methyl ester	382	C25H50O2	002442-49-1	96
4			Tetracosanoic acid, methyl ester	382	C25H50O2	002442-49-1	95
5			Tetracosanoic acid, methyl ester	382	C25H50O2	002442-49-1	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP020325\
Data File : BP023874.D
Acq On : 03 Feb 2025 11:11
Operator : RC/JU
Sample : Q1248-01
Misc :
ALS Vial : 4 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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Methanone, (1-h...	16.345	2.7	ng/ul	715625	4	17.186	5310580	20.0
Hexadecanoic ac...	17.898	3.3	ng/ul	878971	4	17.186	5310580	20.0
n-Hexadecanoic ...	18.128	2.4	ng/ul	638435	4	17.186	5310580	20.0
Hexadecanoic ac...	20.398	3.5	ng/ul	1042740	5	21.639	5998660	20.0
Trifluoroacetox...	21.322	4.2	ng/ul	1258100	5	21.639	5998660	20.0
Docosanoic acid...	21.480	3.6	ng/ul	1069160	5	21.639	5998660	20.0
Tetracosanoic a...	22.769	3.1	ng/ul	924981	5	21.639	5998660	20.0