

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102124\
 Data File : BP022566.D
 Acq On : 23 Oct 2024 16:52
 Operator : RC/JU
 Sample : P4415-06
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 A4F54

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-BP100724.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BP022566.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.193	32	35	42	rVB	16773	21826	0.80%	0.076%
2	3.605	102	105	119	rBV	196006	308351	11.25%	1.068%
3	3.922	155	159	162	rBV3	6418	8457	0.31%	0.029%
4	6.363	570	574	578	rBV7	2308	2929	0.11%	0.010%
5	6.805	643	649	651	rBV2	5623	8543	0.31%	0.030%
6	6.852	651	657	669	rVB	304920	494388	18.03%	1.712%
7	6.999	676	682	694	rVB	216770	346275	12.63%	1.199%
8	7.105	696	700	704	rVB7	1620	3012	0.11%	0.010%
9	7.199	708	716	725	rBV	297074	483756	17.64%	1.675%
10	7.657	786	794	802	rBV	374968	617118	22.51%	2.137%
11	8.357	907	913	923	rBV	380633	658085	24.00%	2.278%
12	8.799	980	988	997	rBV	270500	478868	17.47%	1.658%
13	8.952	1010	1014	1018	rVB7	2278	2945	0.11%	0.010%
14	9.199	1049	1056	1062	rBV3	8893	15960	0.58%	0.055%
15	9.357	1078	1083	1089	rBV4	8722	14681	0.54%	0.051%
16	9.504	1100	1108	1123	rBV	242787	443366	16.17%	1.535%
17	9.769	1149	1153	1160	rVB9	1765	3213	0.12%	0.011%
18	10.046	1191	1200	1211	rBV	429398	727945	26.55%	2.520%
19	10.122	1211	1213	1222	rVB9	5113	10833	0.40%	0.038%
20	10.410	1254	1262	1274	rVB	515593	852412	31.09%	2.951%
21	10.551	1278	1286	1302	rBV	319001	608300	22.19%	2.106%
22	10.957	1346	1355	1365	rBV	4608	13210	0.48%	0.046%
23	11.216	1394	1399	1403	rBV8	1940	4049	0.15%	0.014%
24	11.675	1471	1477	1482	rBV3	5977	9880	0.36%	0.034%
25	11.916	1513	1518	1522	rBV7	1799	3276	0.12%	0.011%
26	12.004	1527	1533	1539	rVB4	5733	10788	0.39%	0.037%
27	12.857	1675	1678	1683	rVB6	2053	3443	0.13%	0.012%
28	13.581	1797	1801	1807	rVB8	3429	5664	0.21%	0.020%
29	13.681	1810	1818	1829	rBV	726070	1151759	42.01%	3.988%
30	13.951	1855	1864	1874	rBV2	810917	1279598	46.67%	4.430%
31	14.163	1896	1900	1906	rVB9	2217	4450	0.16%	0.015%
32	14.263	1910	1917	1924	rBV	879551	1300782	47.44%	4.504%
33	14.328	1924	1928	1934	rVB4	8134	11530	0.42%	0.040%
34	14.392	1934	1939	1940	rBV4	2448	2909	0.11%	0.010%
35	14.504	1951	1958	1980	rBV	253569	463653	16.91%	1.605%
36	14.704	1989	1992	1997	rVB7	7915	12584	0.46%	0.044%

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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_P\Methods\SFAM-EPA-BP100724.MA.M
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37	14.857	2014	2018	2022	rBV7	2050	2918	0.11%	0.010%
38	14.945	2029	2033	2038	rVB8	2316	3836	0.14%	0.013%
39	15.087	2050	2057	2058	rBV6	2318	3756	0.14%	0.013%
40	15.275	2082	2089	2096	rBV	1221089	1739137	63.43%	6.021%
41	15.422	2105	2114	2125	rBV	229444	371272	13.54%	1.285%
42	15.528	2130	2132	2138	rVB5	6590	9485	0.35%	0.033%
43	15.639	2145	2151	2162	rVB	227774	330194	12.04%	1.143%
44	15.786	2172	2176	2179	rBV6	1879	2752	0.10%	0.010%
45	16.022	2211	2216	2221	rVB8	4746	7984	0.29%	0.028%
46	16.069	2222	2224	2230	rVB7	2622	3453	0.13%	0.012%
47	16.169	2238	2241	2245	rBV5	2219	3433	0.13%	0.012%
48	16.469	2289	2292	2298	rBV8	6747	11064	0.40%	0.038%
49	16.657	2320	2324	2329	rVB	15002	19872	0.72%	0.069%
50	16.716	2330	2334	2339	rVB7	4788	9159	0.33%	0.032%
51	16.786	2344	2346	2350	rBV5	2316	4296	0.16%	0.015%
52	16.828	2350	2353	2357	rBV6	3231	6404	0.23%	0.022%
53	16.975	2376	2378	2384	rVB6	4464	6387	0.23%	0.022%
54	17.063	2386	2393	2398	rVV	1115723	1744422	63.62%	6.040%
55	17.104	2398	2400	2404	rVV	75978	95652	3.49%	0.331%
56	17.169	2404	2411	2422	rVB	1249034	2005680	73.15%	6.944%
57	17.257	2423	2426	2427	rBV2	5584	4998	0.18%	0.017%
58	17.281	2427	2430	2435	rVB7	10579	15263	0.56%	0.053%
59	17.451	2456	2459	2462	rBV4	6431	9725	0.35%	0.034%
60	17.675	2493	2497	2499	rBV4	5318	7902	0.29%	0.027%
61	17.880	2527	2532	2537	rBV7	9643	18719	0.68%	0.065%
62	17.998	2545	2552	2562	rBV2	250721	466327	17.01%	1.615%
63	18.169	2577	2581	2585	rVB3	14951	23438	0.85%	0.081%
64	18.586	2647	2652	2662	rVB2	110480	177645	6.48%	0.615%
65	18.745	2676	2679	2684	rBV5	21618	29198	1.06%	0.101%
66	19.098	2734	2739	2743	rBV3	19938	35884	1.31%	0.124%
67	19.204	2748	2757	2762	rBV	193314	300488	10.96%	1.040%
68	19.263	2762	2767	2773	rVV	52822	92758	3.38%	0.321%
69	19.327	2775	2778	2789	rVV2	118536	181600	6.62%	0.629%
70	19.551	2810	2816	2825	rBV	1985921	2741763	100.00%	9.493%
71	20.022	2891	2896	2900	rBV	144927	198195	7.23%	0.686%
72	21.174	3089	3092	3100	rVB4	107027	173621	6.33%	0.601%
73	21.504	3140	3148	3153	rBV	1270489	2216568	80.84%	7.674%
74	21.804	3196	3199	3203	rVB	85890	112165	4.09%	0.388%
75	22.357	3289	3293	3306	rVB2	140155	276781	10.10%	0.958%
76	23.898	3550	3555	3563	rVB2	162924	362902	13.24%	1.256%

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77	24.533	3655	3663	3673	rBV	1008457	2513342	91.67%	8.702%
78	24.768	3694	3703	3712	rVB	760010	2154174	78.57%	7.458%

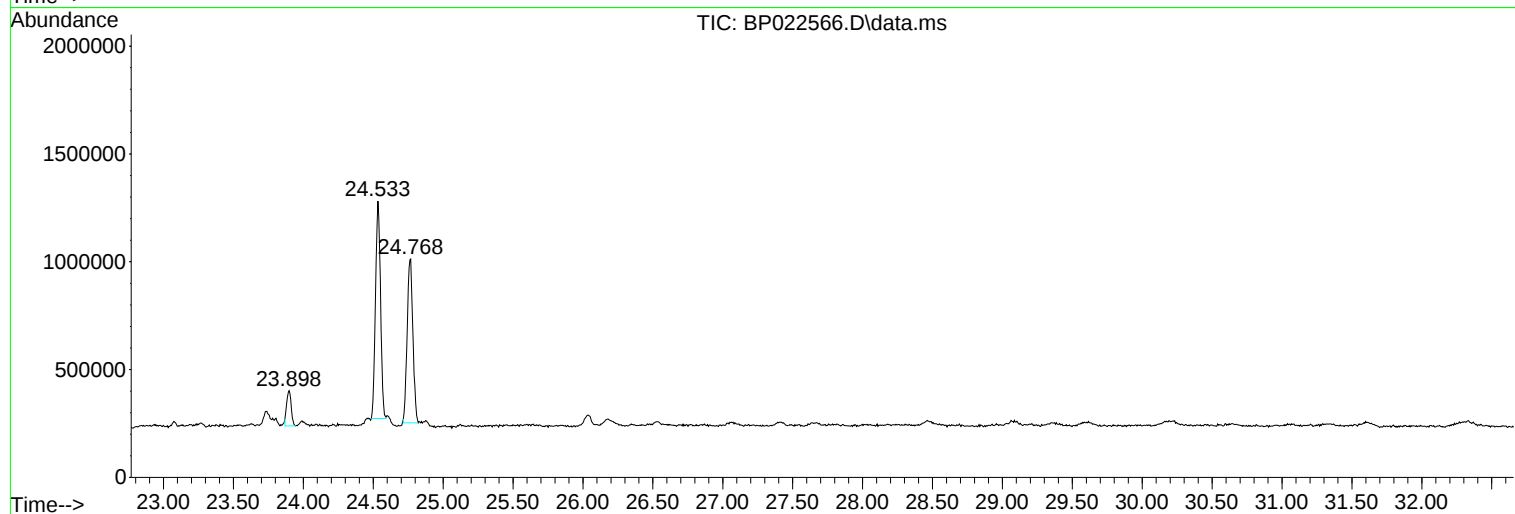
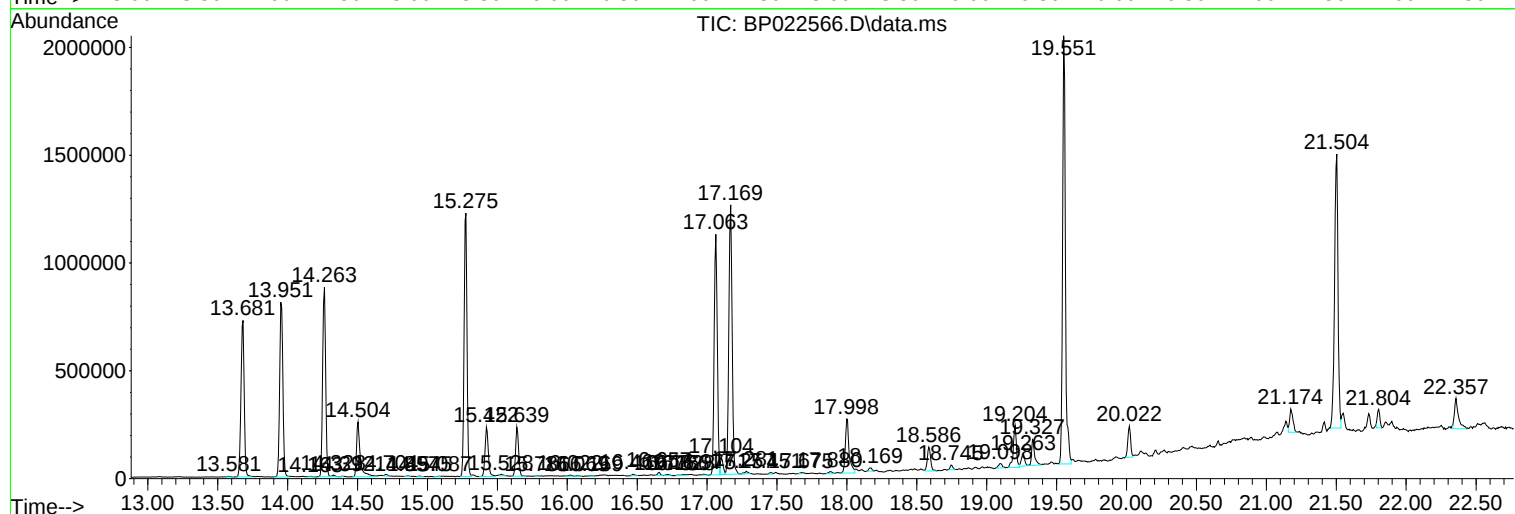
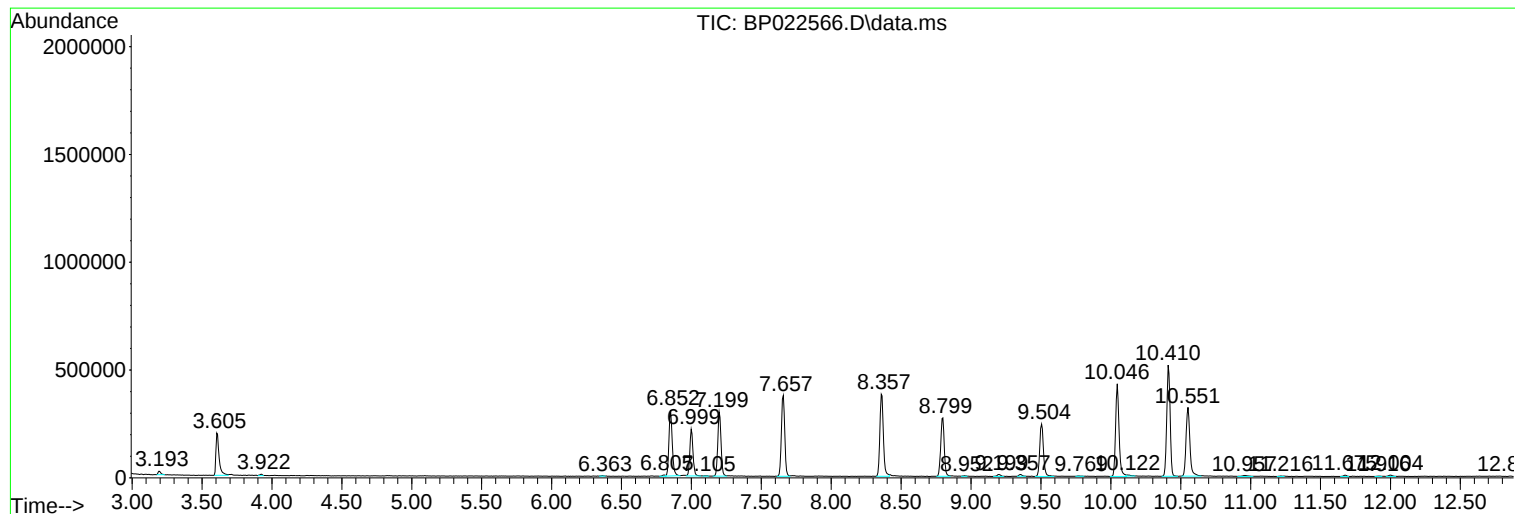
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100724.MA.M
Quant Title : SVOA CALIBRATION

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



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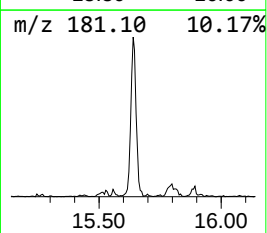
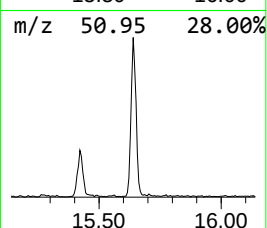
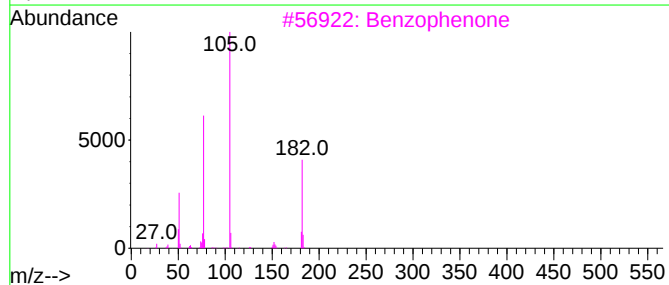
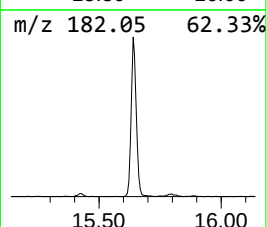
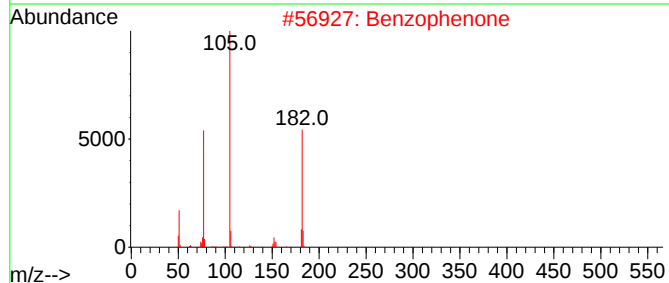
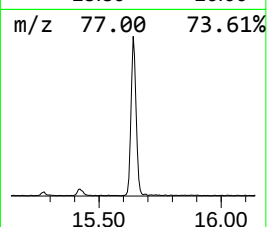
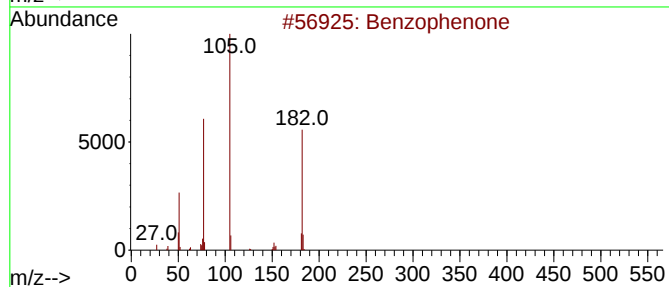
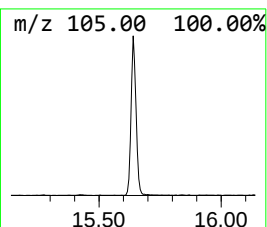
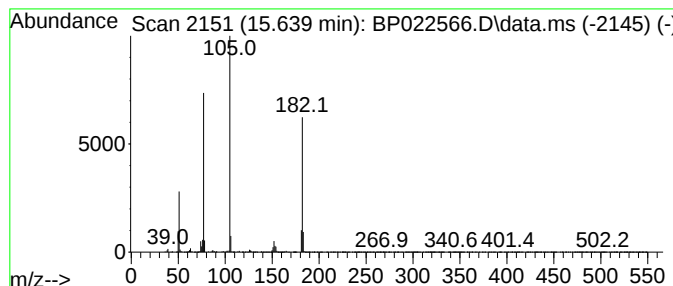
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100724.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Benzophenone Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.639	5.08 ng/ul	330194	Acenaphthene-d10	14.263

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzophenone	182	C13H10O	000119-61-9	96
2			Benzophenone	182	C13H10O	000119-61-9	95
3			Benzophenone	182	C13H10O	000119-61-9	94
4			Benzophenone	182	C13H10O	000119-61-9	94
5			Benzophenone	182	C13H10O	000119-61-9	91



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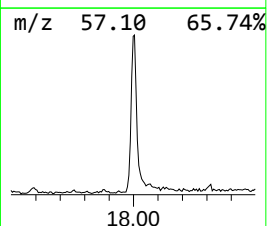
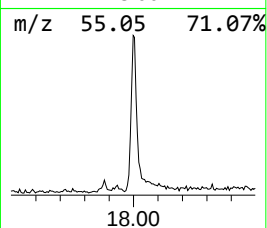
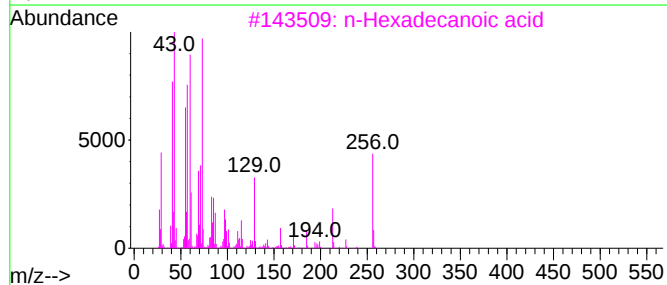
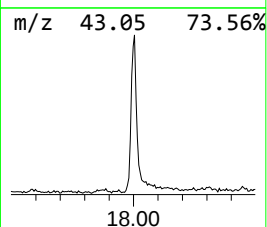
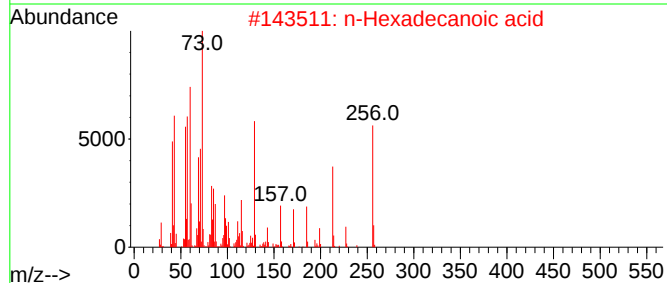
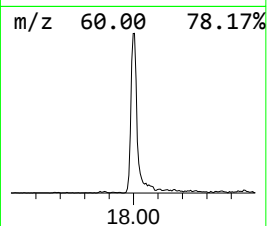
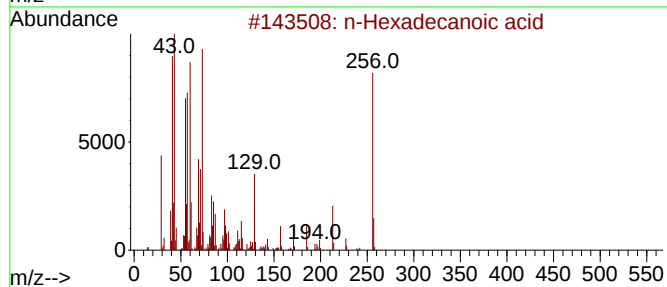
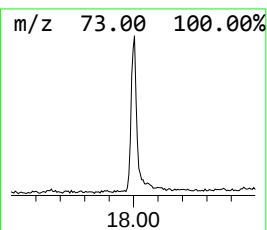
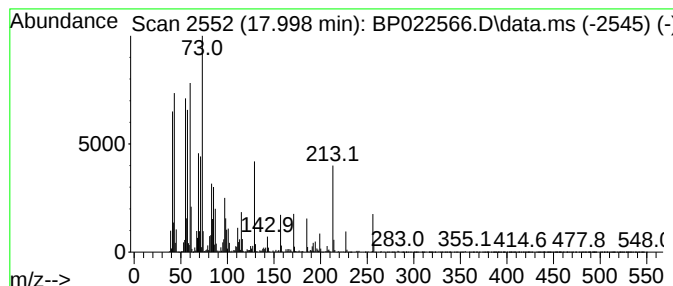
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100724.MA.M
Quant Title : SVOA CALIBRATION

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.
17.998	5.35 ng/ul	466327	Phenanthrene-d10	17.063

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	97		
3	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	96		
4	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	95		
5	Tridecanoic acid	214	C13H26O2	000638-53-9	93		



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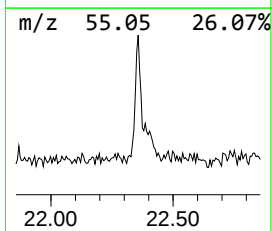
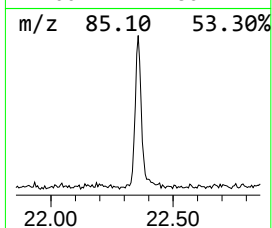
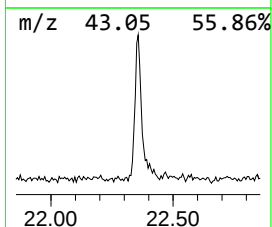
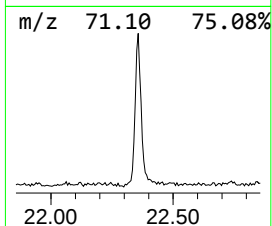
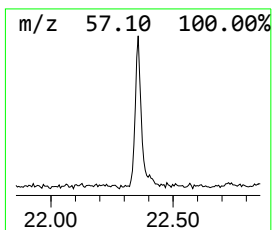
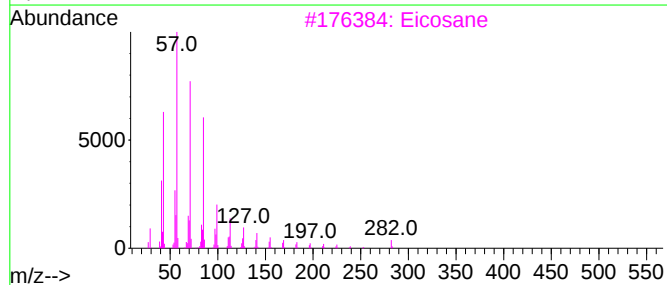
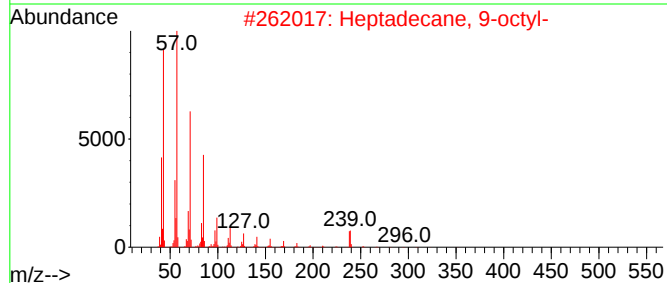
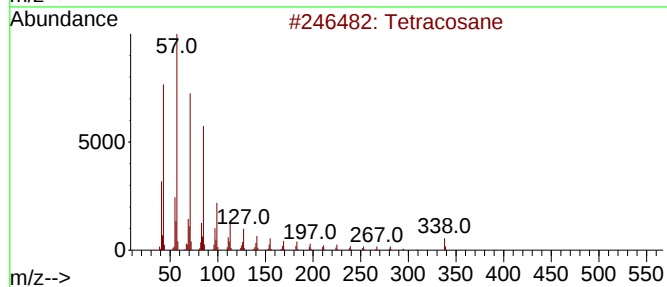
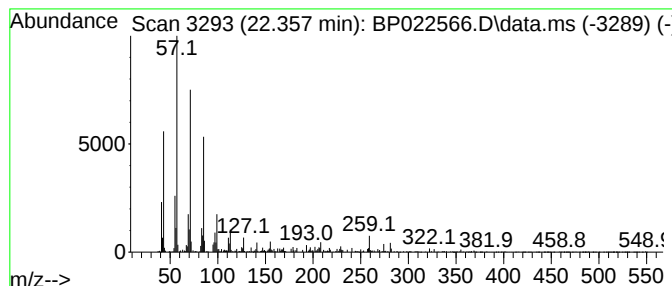
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100724.MA.M
Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.357	2.50 ng/ul	276781	Chrysene-d12	21.504

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tetracosane	338	C24H50	000646-31-1	95
2			Heptadecane, 9-octyl-	352	C25H52	007225-64-1	93
3			Eicosane	282	C20H42	000112-95-8	93
4			2-Methyltetracosane	352	C25H52	001560-78-7	93
5			Nonadecane	268	C19H40	000629-92-5	93



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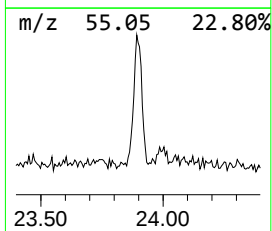
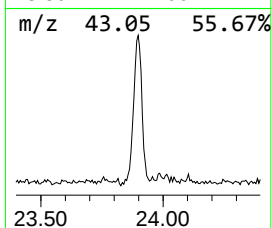
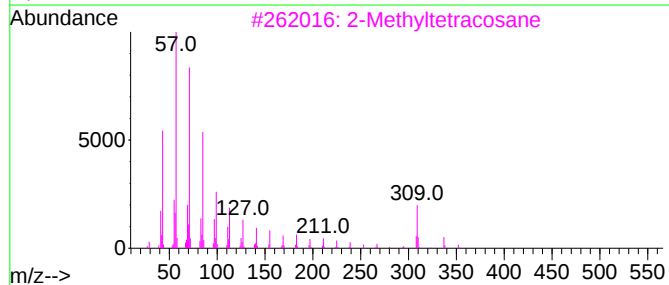
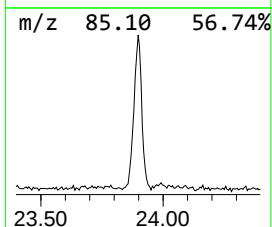
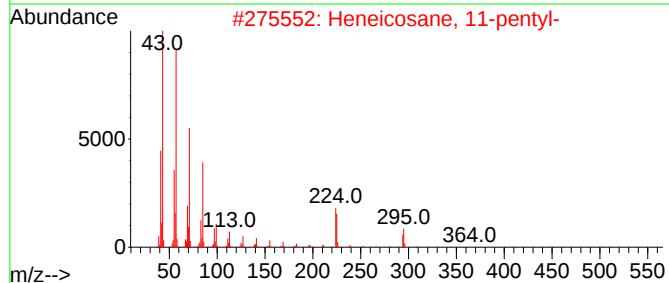
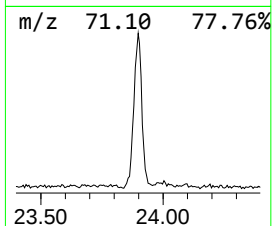
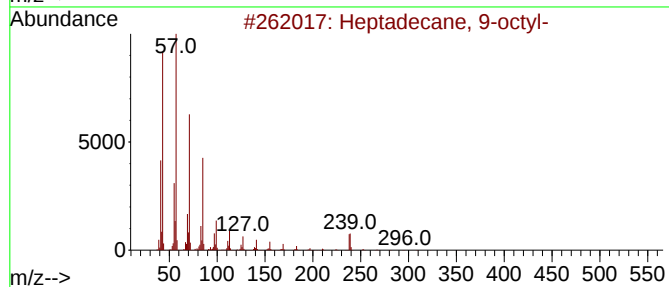
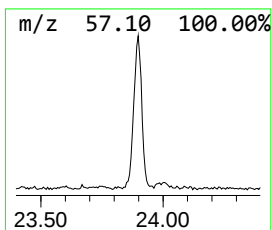
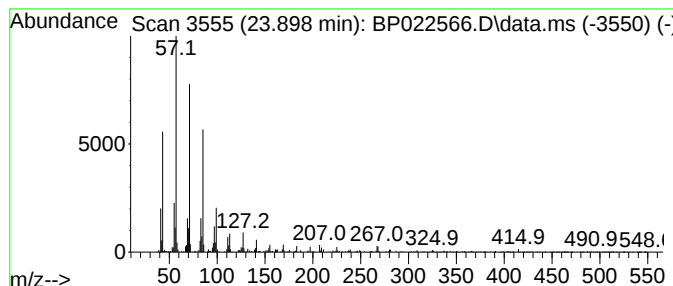
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100724.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 5 (DEL) Alkane: Straight-Chai... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.898	3.37 ng/ul	362902	Perylene-d12	24.768

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heptadecane, 9-octyl-	352	C25H52	007225-64-1	93
2			Heneicosane, 11-pentyl-	366	C26H54	014739-72-1	93
3			2-Methyltetracosane	352	C25H52	001560-78-7	91
4			Heneicosane	296	C21H44	000629-94-7	91
5			Hexacosane	366	C26H54	000630-01-3	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102124\
Data File : BP022566.D
Acq On : 23 Oct 2024 16:52
Operator : RC/JU
Sample : P4415-06
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A4F54

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100724.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
Benzophenone	15.639	5.1	ng/u1	330194	3	14.263	1300780	20.0
n-Hexadecanoic ...	17.998	5.3	ng/u1	466327	4	17.063	1744420	20.0
(DEL) Alkane: S...	22.357	2.5	ng/u1	276781	5	21.504	2216570	20.0
(DEL) Alkane: S...	23.898	3.4	ng/u1	362902	6	24.768	2154170	20.0