

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM102824\
 Data File : BM048285.D
 Acq On : 29 Oct 2024 12:47
 Operator : RC/JU
 Sample : P4532-16
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 GCNX7

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-BM102624.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BM048285.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.169	27	31	35	rBV	61020	73430	1.38%	0.112%
2	3.587	99	102	112	rBV	261845	405867	7.64%	0.622%
3	3.910	154	157	164	rVB	12201	17016	0.32%	0.026%
4	4.287	216	221	227	rBV6	12292	20464	0.39%	0.031%
5	4.469	249	252	256	rVV3	6381	9559	0.18%	0.015%
6	5.428	407	415	419	rBV9	5387	12412	0.23%	0.019%
7	5.522	427	431	436	rBV4	15822	24510	0.46%	0.038%
8	5.740	460	468	474	rBV4	7521	13952	0.26%	0.021%
9	5.851	483	487	491	rVB7	6889	11767	0.22%	0.018%
10	6.069	518	524	528	rBV	49587	78044	1.47%	0.120%
11	6.293	557	562	563	rBV4	8668	11615	0.22%	0.018%
12	6.528	596	602	606	rVV2	23120	37274	0.70%	0.057%
13	6.575	606	610	617	rVB2	30301	49425	0.93%	0.076%
14	6.745	634	639	647	rBV2	4036	11004	0.21%	0.017%
15	6.892	656	664	676	rBV	241379	436937	8.22%	0.669%
16	7.051	683	691	703	rBV	761196	1230039	23.14%	1.884%
17	7.251	718	725	737	rBV	842042	1382513	26.01%	2.118%
18	7.334	737	739	744	rVV6	7035	12774	0.24%	0.020%
19	7.716	795	804	812	rBV	754459	1231242	23.17%	1.886%
20	8.428	919	925	943	rVV	728061	1249050	23.50%	1.913%
21	8.875	994	1001	1009	rBV	864109	1445438	27.20%	2.214%
22	8.945	1009	1013	1018	rVB	55075	81148	1.53%	0.124%
23	9.128	1018	1044	1062	rBV	1240509	4992481	93.94%	7.647%
24	9.298	1069	1073	1078	rVB	27938	44548	0.84%	0.068%
25	9.445	1093	1098	1107	rVB6	19869	44828	0.84%	0.069%
26	9.592	1116	1123	1142	rBV	732002	1261487	23.74%	1.932%
27	9.839	1148	1165	1179	rBV	425141	1353789	25.47%	2.074%
28	9.933	1179	1181	1191	rVB5	16276	29270	0.55%	0.045%
29	10.039	1196	1199	1205	rVB5	10270	13021	0.25%	0.020%
30	10.133	1207	1215	1227	rBV	1093060	1961481	36.91%	3.004%
31	10.504	1269	1278	1289	rBV	1115916	1989694	37.44%	3.048%
32	10.645	1295	1302	1322	rBV	819892	1605042	30.20%	2.458%
33	10.798	1322	1328	1336	rVB10	13589	29518	0.56%	0.045%
34	11.345	1412	1421	1427	rBV10	13520	30221	0.57%	0.046%
35	11.463	1430	1441	1451	rBV	302726	742186	13.97%	1.137%
36	11.645	1469	1472	1481	rVB8	7835	16850	0.32%	0.026%

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Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM102624.MA.M
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37	11.792	1493	1497	1502	rBV3	11466	20340	0.38%	0.031%
38	11.933	1515	1521	1526	rBV5	9206	15025	0.28%	0.023%
39	12.033	1533	1538	1540	rBV6	8055	12240	0.23%	0.019%
40	12.104	1543	1550	1556	rVB	18922	33531	0.63%	0.051%

41	12.263	1572	1577	1581	rBV3	11155	18991	0.36%	0.029%
42	12.351	1587	1592	1598	rVB7	5095	8401	0.16%	0.013%
43	12.927	1684	1690	1694	rVB2	8042	13610	0.26%	0.021%
44	13.263	1739	1747	1755	rBV	167998	269238	5.07%	0.412%
45	13.774	1827	1834	1842	rBV	2037131	3006982	56.58%	4.606%

46	13.880	1849	1852	1857	rVB6	5037	8321	0.16%	0.013%
47	14.051	1872	1881	1892	rBV2	2411739	3541680	66.64%	5.425%
48	14.204	1896	1907	1924	rVV	1294576	2611497	49.14%	4.000%
49	14.357	1926	1933	1944	rVB	1666500	2489609	46.85%	3.813%
50	14.486	1950	1955	1961	rVB9	9109	16803	0.32%	0.026%

51	14.592	1967	1973	1983	rBV	129815	322702	6.07%	0.494%
52	15.021	2042	2046	2049	rBV5	7335	9137	0.17%	0.014%
53	15.092	2054	2058	2066	rBV3	16356	39482	0.74%	0.060%
54	15.163	2066	2070	2073	rVV5	8646	14375	0.27%	0.022%
55	15.351	2093	2102	2118	rVV	3152784	4497303	84.62%	6.889%

56	15.480	2118	2124	2140	rVV	1024215	1479883	27.85%	2.267%
57	15.592	2140	2143	2151	rVB4	18835	29336	0.55%	0.045%
58	15.715	2158	2164	2173	rBV	457694	656621	12.36%	1.006%
59	15.868	2186	2190	2194	rBV2	21551	32092	0.60%	0.049%
60	15.915	2195	2198	2202	rVB3	13137	18731	0.35%	0.029%

61	16.057	2218	2222	2228	rBV3	19852	36413	0.69%	0.056%
62	16.139	2235	2236	2241	rVB5	9179	10281	0.19%	0.016%
63	16.280	2253	2260	2267	rBV	98043	139245	2.62%	0.213%
64	16.380	2273	2277	2279	rBV4	12496	17245	0.32%	0.026%
65	16.421	2281	2284	2288	rVB5	12496	17643	0.33%	0.027%

66	16.515	2296	2300	2304	rBV6	9612	16174	0.30%	0.025%
67	16.568	2304	2309	2321	rBV	187769	295903	5.57%	0.453%
68	16.704	2326	2332	2340	rBV2	85348	143826	2.71%	0.220%
69	16.868	2354	2360	2363	rBV3	25625	33414	0.63%	0.051%
70	16.921	2365	2369	2373	rVB7	12477	20982	0.39%	0.032%

71	16.980	2377	2379	2382	rVB4	9057	8147	0.15%	0.012%
72	17.104	2393	2400	2409	rBV	1827479	2641753	49.71%	4.046%
73	17.198	2410	2416	2428	rBV	3228032	4741018	89.21%	7.262%
74	17.386	2444	2448	2454	rVB3	18832	32694	0.62%	0.050%
75	17.527	2468	2472	2476	rBV4	25175	34211	0.64%	0.052%

76	17.574	2476	2480	2484	rBV2	20481	32452	0.61%	0.050%
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 Title : SVOA CALIBRATION

77	17.668	2493	2496	2499	rBV4	13760	22993	0.43%	0.035%
78	17.774	2511	2514	2518	rBV4	15536	21856	0.41%	0.033%
79	17.815	2518	2521	2526	rVB7	15152	25426	0.48%	0.039%
80	17.898	2533	2535	2539	rVB5	9406	10190	0.19%	0.016%
81	17.945	2539	2543	2544	rBV3	6584	8628	0.16%	0.013%
82	17.998	2547	2552	2555	rBV	94790	137530	2.59%	0.211%
83	18.039	2555	2559	2563	rVB	85247	117541	2.21%	0.180%
84	18.280	2596	2600	2612	rVB2	60354	119452	2.25%	0.183%
85	18.533	2639	2643	2647	rBV4	24105	32664	0.61%	0.050%
86	18.603	2653	2655	2659	rVB5	8175	8501	0.16%	0.013%
87	18.662	2662	2665	2667	rVV3	12804	13900	0.26%	0.021%
88	18.845	2692	2696	2702	rVB4	27253	46550	0.88%	0.071%
89	18.986	2717	2720	2727	rVB	29836	46186	0.87%	0.071%
90	19.056	2728	2732	2741	rBV2	40343	74356	1.40%	0.114%
91	19.133	2741	2745	2746	rBV	46045	56825	1.07%	0.087%
92	19.156	2746	2749	2754	rVB	69724	85927	1.62%	0.132%
93	19.280	2766	2770	2778	rBV6	44767	65015	1.22%	0.100%
94	19.427	2792	2795	2800	rVB3	30207	37179	0.70%	0.057%
95	19.492	2800	2806	2818	rBV	3874206	5314514	100.00%	8.140%
96	19.603	2822	2825	2828	rBV4	19890	26361	0.50%	0.040%
97	21.015	3061	3065	3071	rBV4	75649	132642	2.50%	0.203%
98	21.327	3112	3118	3132	rBV	1614694	2488592	46.83%	3.812%
99	24.068	3576	3584	3600	rVB	1862200	4499612	84.67%	6.892%
100	24.268	3610	3618	3636	rVB	977085	2545946	47.91%	3.900%

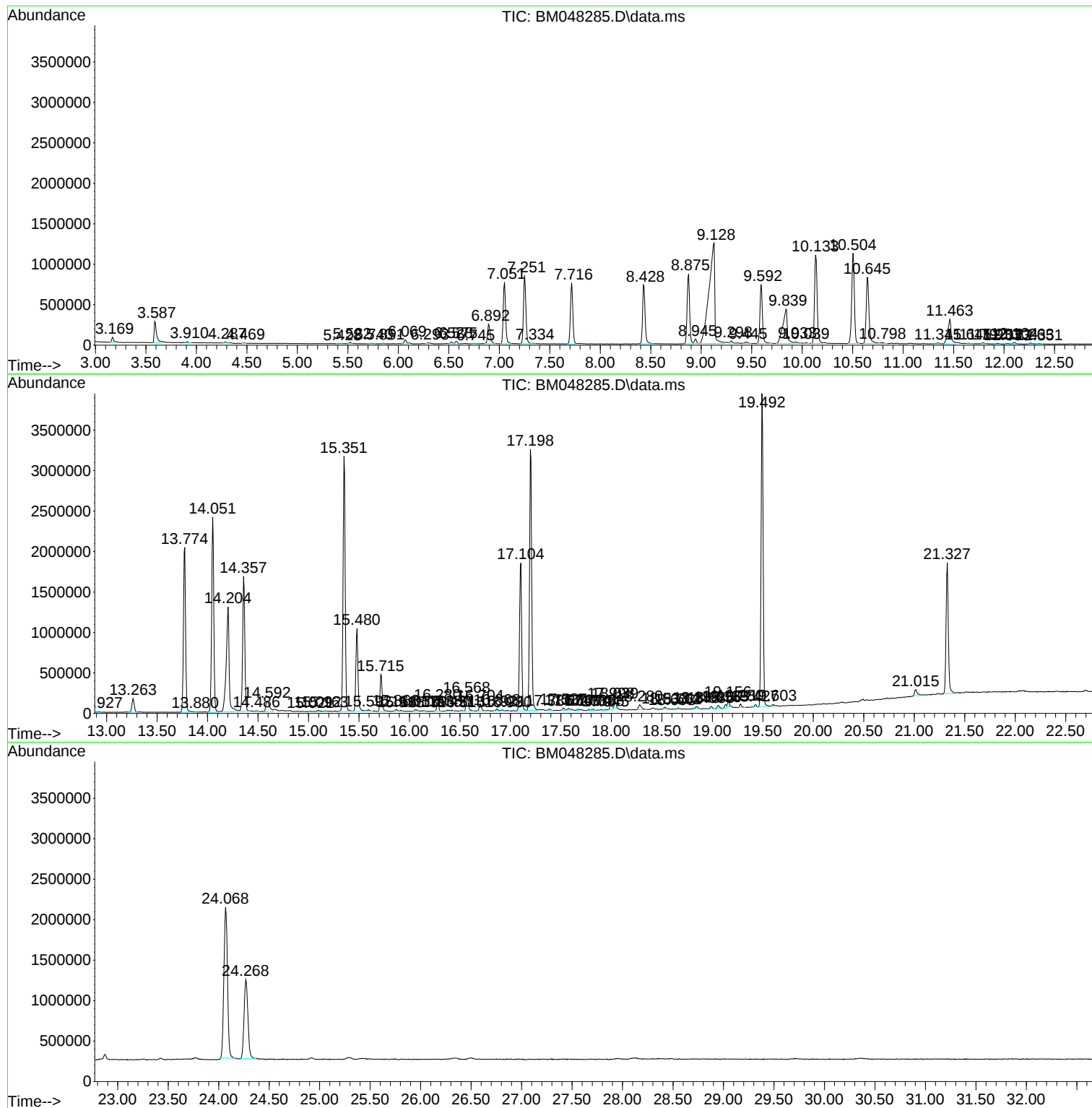
Sum of corrected areas: 65285613

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM102624.MA.M
Quant Title : SVOA CALIBRATION

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



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ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_M
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GCNX7

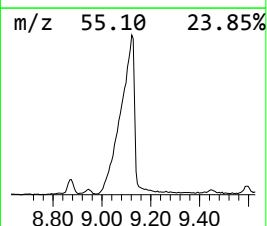
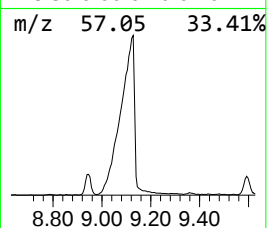
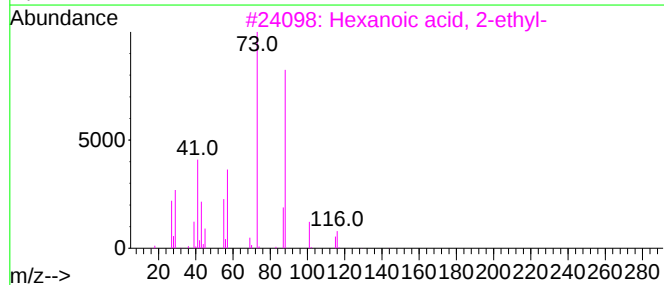
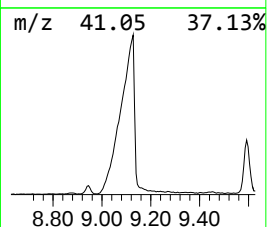
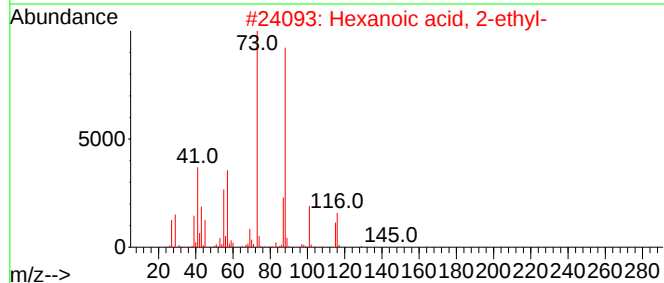
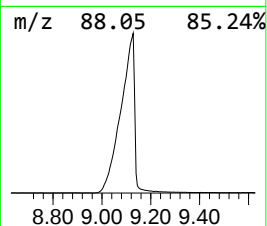
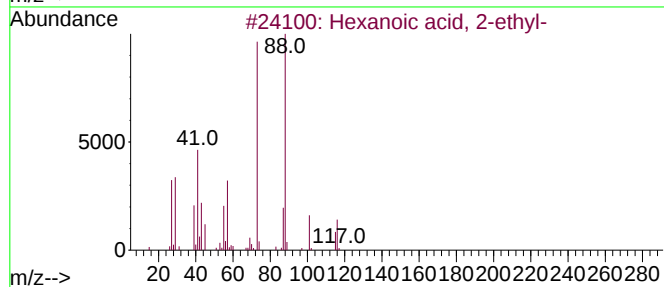
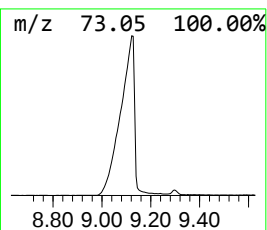
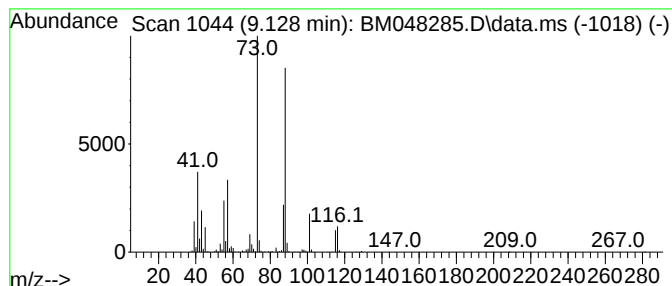
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM102624.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Hexanoic acid, 2-ethyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.128	50.18 ng/ul	4992480	Naphthalene-d8	10.504

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexanoic acid, 2-ethyl-	144	C8H16O2	000149-57-5	90
2			Hexanoic acid, 2-ethyl-	144	C8H16O2	000149-57-5	90
3			Hexanoic acid, 2-ethyl-	144	C8H16O2	000149-57-5	90
4			Hexanoic acid, 2-ethyl-	144	C8H16O2	000149-57-5	86
5			Hexanoic acid, 2-ethyl-	144	C8H16O2	000149-57-5	86



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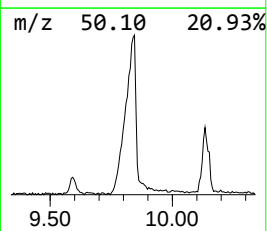
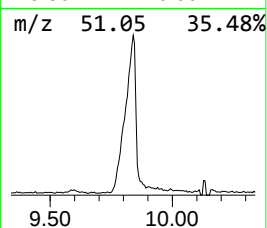
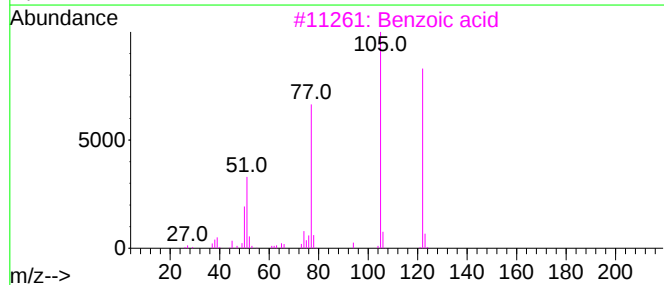
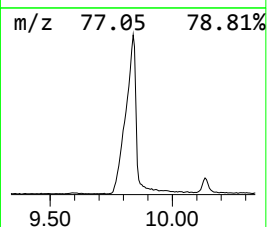
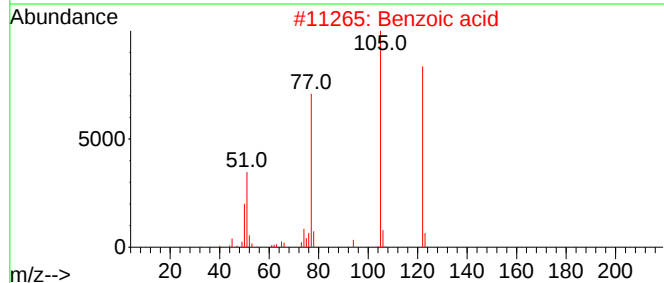
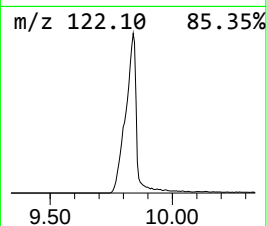
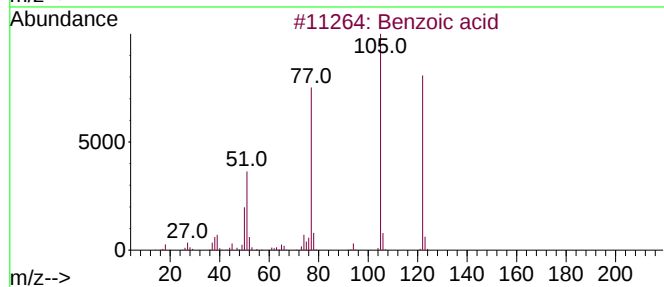
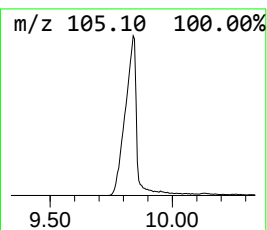
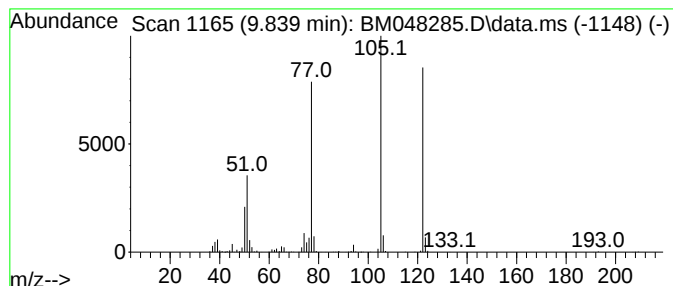
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM102624.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 Benzoic acid Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.839	13.61 ng/ul	1353790	Naphthalene-d8	10.504

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzoic acid	122	C7H6O2	000065-85-0	96
2			Benzoic acid	122	C7H6O2	000065-85-0	96
3			Benzoic acid	122	C7H6O2	000065-85-0	96
4			Benzoic acid	122	C7H6O2	000065-85-0	94
5			Benzoic acid, silver(1+) salt	228	C7H5AgO2	000532-31-0	90



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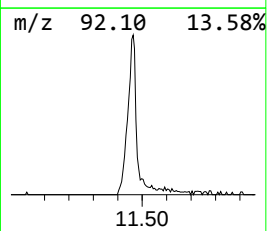
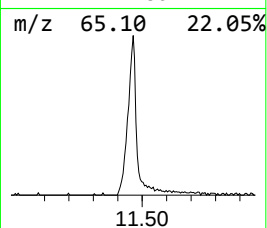
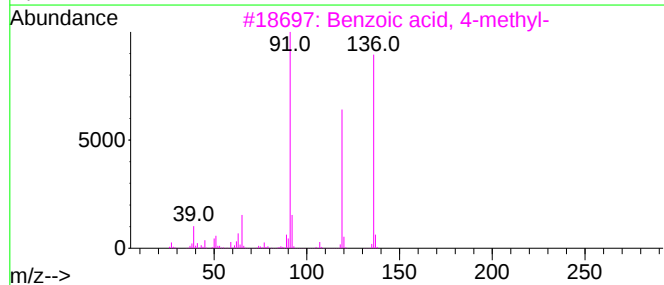
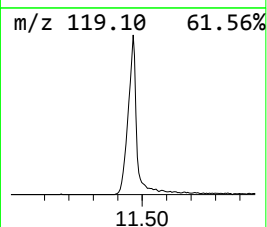
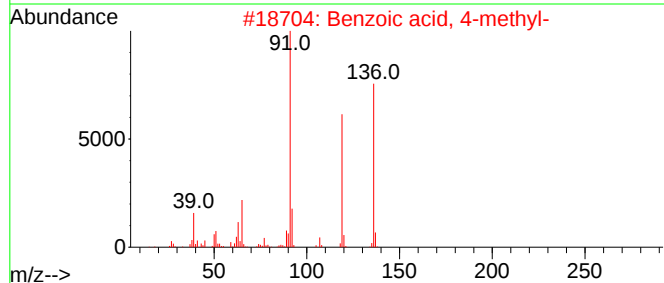
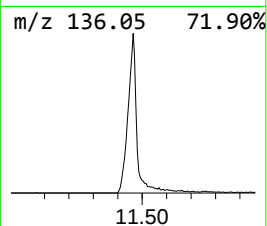
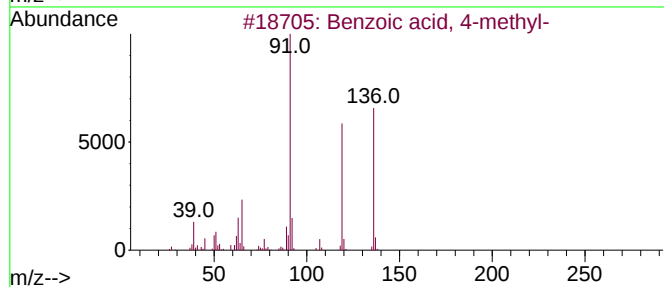
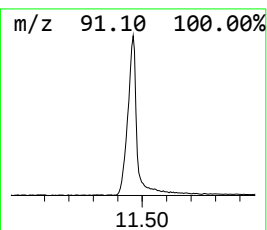
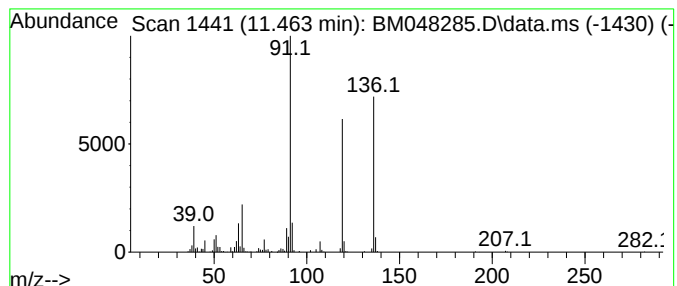
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Quant Title : SVOA CALIBRATION

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

Peak Number 3 Benzoic acid, 4-methyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.463	7.46 ng/ul	742186	Naphthalene-d8	10.504

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzoic acid, 4-methyl-	136	C8H8O2	000099-94-5	96
2			Benzoic acid, 4-methyl-	136	C8H8O2	000099-94-5	96
3			Benzoic acid, 4-methyl-	136	C8H8O2	000099-94-5	95
4			Benzoic acid, 4-methyl-	136	C8H8O2	000099-94-5	94
5			Benzoic acid, 4-methyl-	136	C8H8O2	000099-94-5	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM102824\
Data File : BM048285.D
Acq On : 29 Oct 2024 12:47
Operator : RC/JU
Sample : P4532-16
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
GCNX7

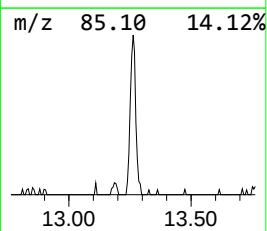
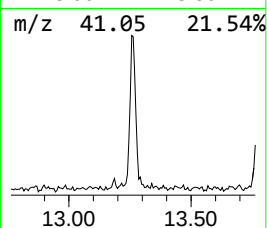
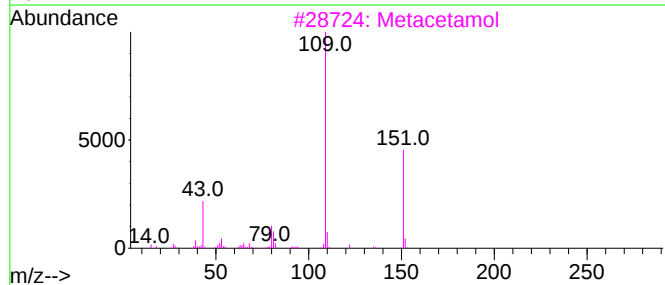
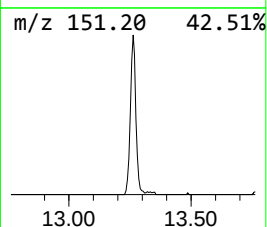
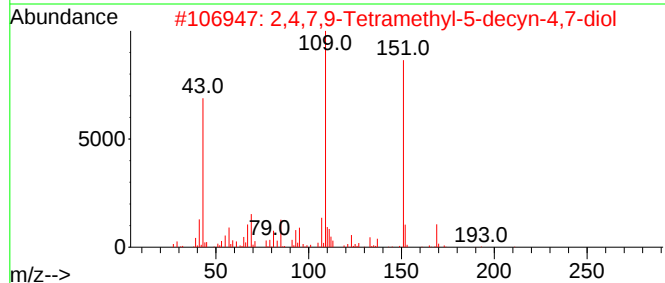
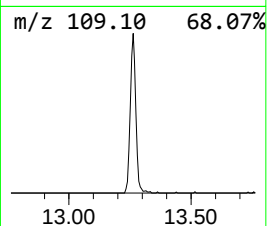
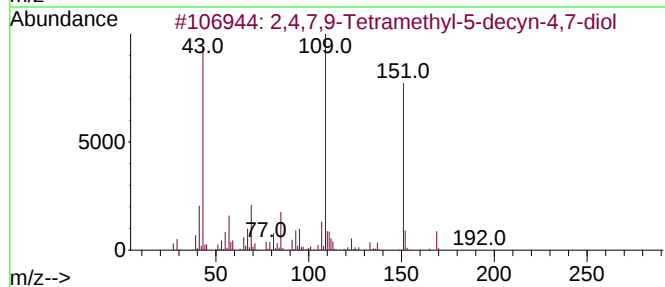
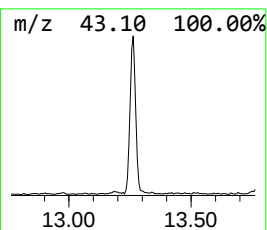
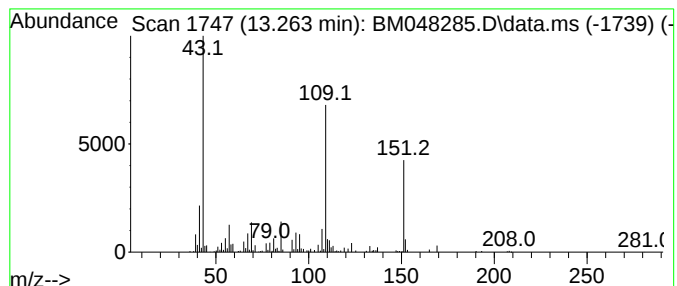
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 2,4,7,9-Tetramethyl-5-decyn... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.263	2.16 ng/ul	269238	Acenaphthene-d10	14.357

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,4,7,9-Tetramethyl-5-decyn-4,7-...	226	C14H26O2	000126-86-3	91		
2	2,4,7,9-Tetramethyl-5-decyn-4,7-...	226	C14H26O2	000126-86-3	59		
3	Metacetamol	151	C8H9NO2	000621-42-1	52		
4	Metacetamol	151	C8H9NO2	000621-42-1	50		
5	Acetaminophen	151	C8H9NO2	000103-90-2	50		



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM102824\
Data File : BM048285.D
Acq On : 29 Oct 2024 12:47
Operator : RC/JU
Sample : P4532-16
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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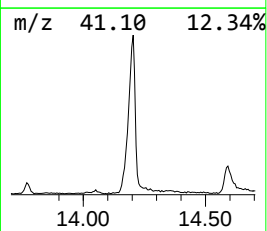
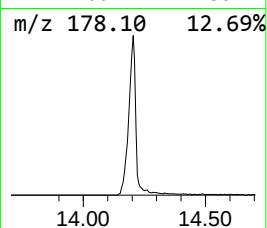
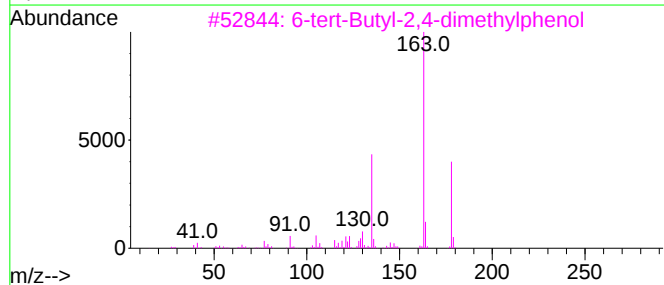
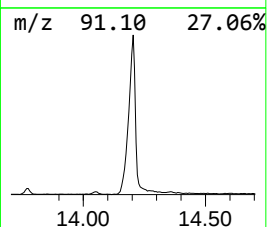
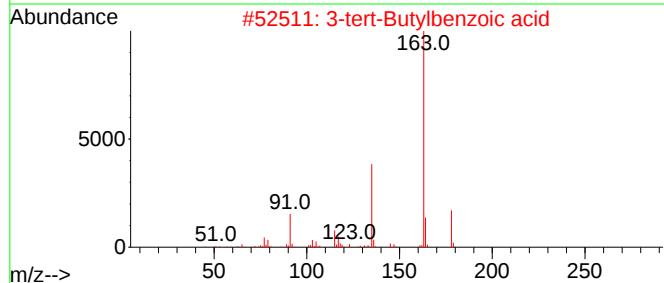
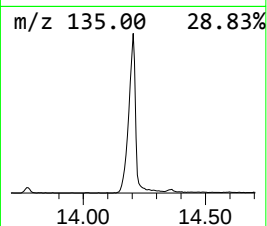
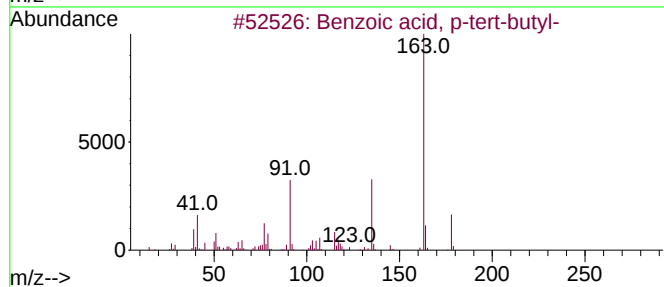
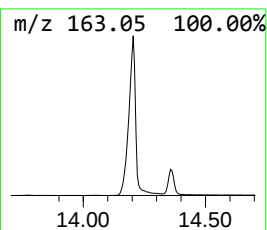
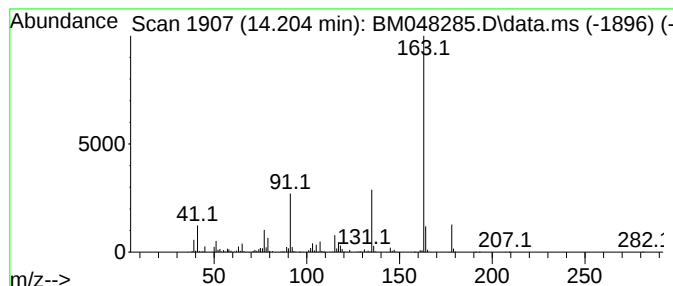
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 Benzoic acid, p-tert-butyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.204	20.98 ng/ul	2611500	Acenaphthene-d10	14.357

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzoic acid, p-tert-butyl-	178	C11H14O2	000098-73-7	96
2			3-tert-Butylbenzoic acid	178	C11H14O2	007498-54-6	94
3			6-tert-Butyl-2,4-dimethylphenol	178	C12H18O	001879-09-0	81
4			Benzoic acid, 4-acetyl-, methyl ...	178	C10H10O3	003609-53-8	72
5			Benzaldehyde, 5-(1,1-dimethyleth...	178	C11H14O2	002725-53-3	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM102824\
Data File : BM048285.D
Acq On : 29 Oct 2024 12:47
Operator : RC/JU
Sample : P4532-16
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
GCNX7

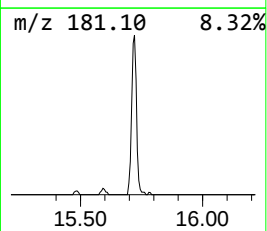
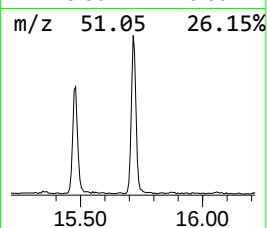
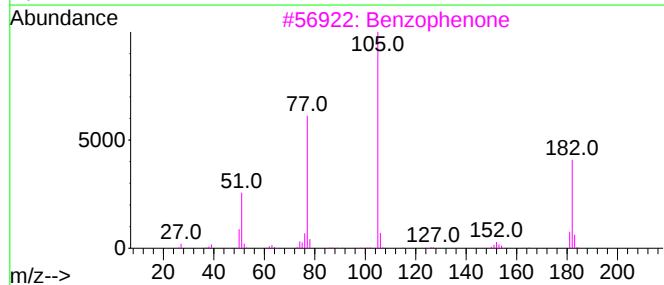
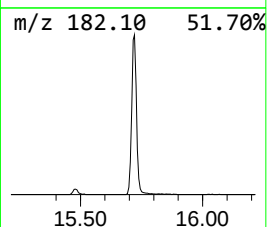
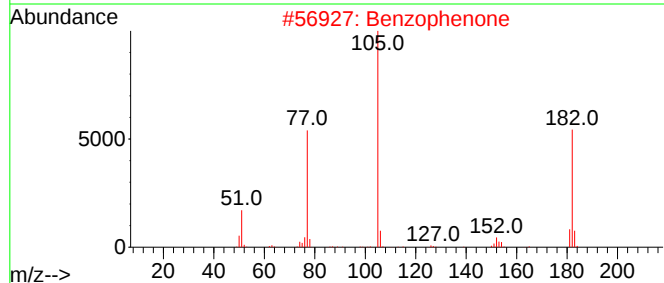
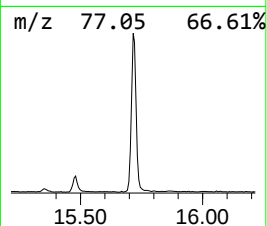
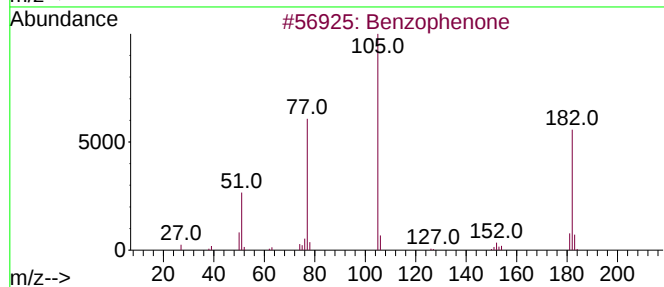
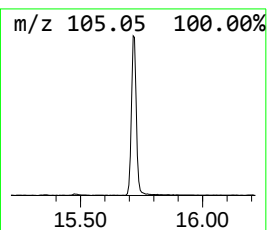
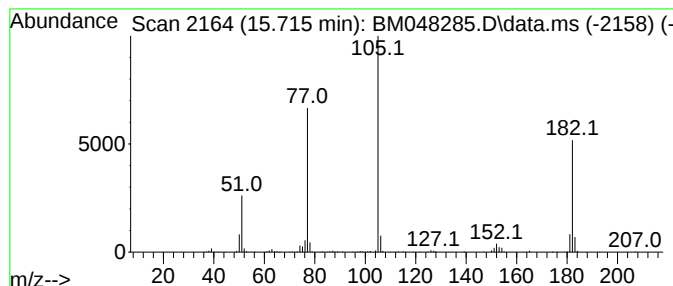
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 Benzophenone Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.715	5.27 ng/ul	656621	Acenaphthene-d10	14.357

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	90
5			Azobenzene	182	C12H10N2	000103-33-3	72



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM102824\
Data File : BM048285.D
Acq On : 29 Oct 2024 12:47
Operator : RC/JU
Sample : P4532-16
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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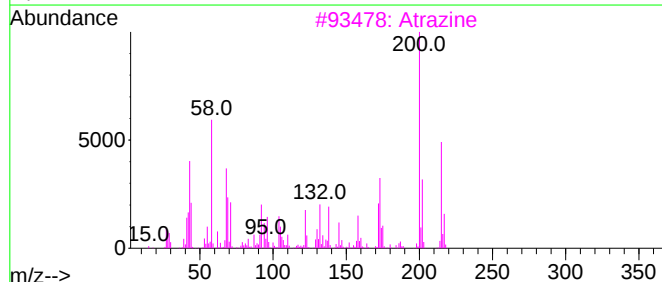
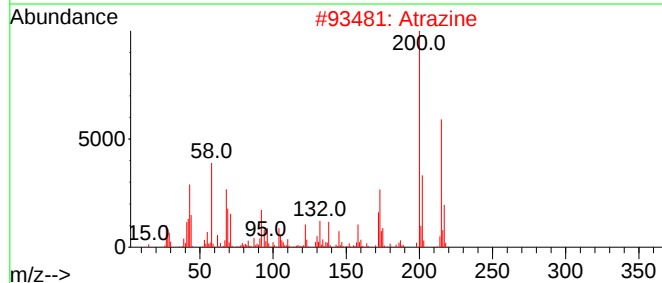
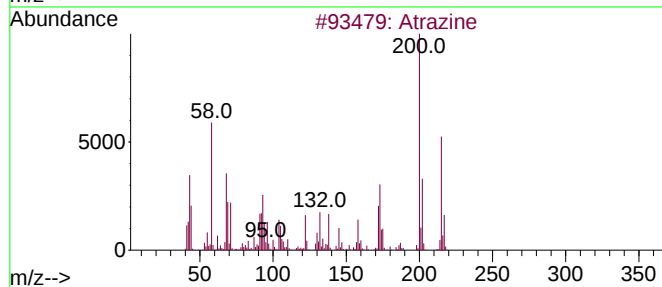
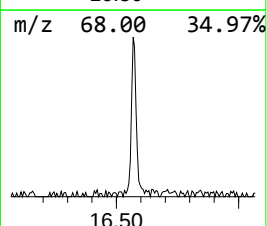
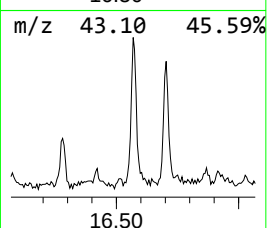
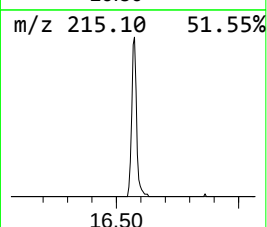
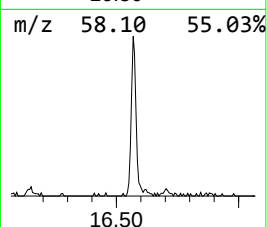
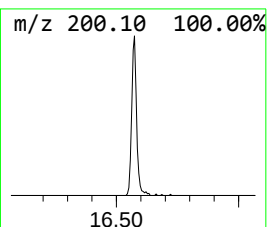
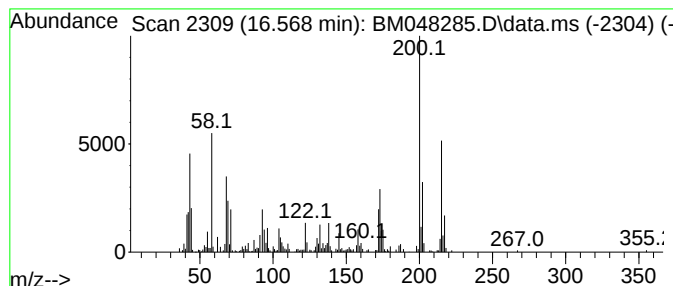
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 Atrazine Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.568	2.24 ng/ul	295903	Phenanthrene-d10	17.104

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Atrazine			215	C8H14ClN5	001912-24-9	99
2	Atrazine			215	C8H14ClN5	001912-24-9	99
3	Atrazine			215	C8H14ClN5	001912-24-9	99
4	Atrazine			215	C8H14ClN5	001912-24-9	98
5	Atrazine			215	C8H14ClN5	001912-24-9	98



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM102824\
Data File : BM048285.D
Acq On : 29 Oct 2024 12:47
Operator : RC/JU
Sample : P4532-16
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
GCNX7

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM102624.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
Hexanoic acid, ...	9.128	50.2	ng/ul	4992480	2	10.504	1989690	20.0
Benzoic acid	9.839	13.6	ng/ul	1353790	2	10.504	1989690	20.0
Benzoic acid, 4...	11.463	7.5	ng/ul	742186	2	10.504	1989690	20.0
2,4,7,9-Tetrame...	13.263	2.2	ng/ul	269238	3	14.357	2489610	20.0
Benzoic acid, p...	14.204	21.0	ng/ul	2611500	3	14.357	2489610	20.0
Benzophenone	15.715	5.3	ng/ul	656621	3	14.357	2489610	20.0
Atra zine	16.568	2.2	ng/ul	295903	4	17.104	2641750	20.0