

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM020322\
 Data File : BM034011.D
 Acq On : 03 Feb 2022 20:34
 Operator : CG/JU
 Sample : N1365-02
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 BGLY8

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-BM013122.M
 Title : SVOA CALIBRATION

Signal : TIC: BM034011.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.284	5	8	15	rVB	19797	24804	1.15%	0.091%
2	3.719	79	82	95	rBV	105470	181309	8.41%	0.668%
3	3.825	96	100	108	rVB	27064	39642	1.84%	0.146%
4	3.961	118	123	126	rBV3	10103	16065	0.74%	0.059%
5	4.055	135	139	144	rVB	21178	28468	1.32%	0.105%
6	4.225	164	168	173	rBV5	5392	8835	0.41%	0.033%
7	4.284	174	178	182	rVB3	6714	9106	0.42%	0.034%
8	4.596	228	231	239	rVB8	6718	15392	0.71%	0.057%
9	4.996	294	299	304	rVB3	20485	32055	1.49%	0.118%
10	5.090	310	315	320	rBV7	5867	11308	0.52%	0.042%
11	6.919	613	626	630	rBV9	4434	13665	0.63%	0.050%
12	7.078	646	653	664	rVB	340583	577912	26.79%	2.128%
13	7.243	674	681	690	rBV	265237	411003	19.05%	1.514%
14	7.449	709	716	725	rBV	346485	549440	25.47%	2.024%
15	7.543	728	732	736	rBV5	6702	11877	0.55%	0.044%
16	7.919	789	796	803	rBV	504419	780418	36.18%	2.874%
17	7.984	803	807	814	rVB	26077	39935	1.85%	0.147%
18	8.625	908	916	925	rBV	405430	661107	30.65%	2.435%
19	8.743	932	936	946	rVB	17221	29350	1.36%	0.108%
20	9.078	986	993	1001	rBV	319879	531432	24.64%	1.957%
21	9.201	1010	1014	1018	rVB4	7531	10673	0.49%	0.039%
22	9.531	1063	1070	1075	rBV2	19858	32069	1.49%	0.118%
23	9.807	1109	1117	1123	rBV	277026	444589	20.61%	1.637%
24	10.348	1202	1209	1217	rBV	434176	718674	33.32%	2.647%
25	10.507	1231	1236	1245	rBV	56662	101325	4.70%	0.373%
26	10.731	1266	1274	1280	rBV	669744	1118267	51.84%	4.119%
27	10.778	1280	1282	1289	rVB2	12200	17577	0.81%	0.065%
28	10.860	1289	1296	1311	rBV	300482	573789	26.60%	2.113%
29	11.742	1441	1446	1456	rVB	25592	46490	2.16%	0.171%
30	12.160	1512	1517	1523	rVV	9696	14787	0.69%	0.054%
31	12.319	1539	1544	1549	rVV	8605	13058	0.61%	0.048%
32	12.389	1552	1556	1562	rVB2	10755	15320	0.71%	0.056%
33	12.495	1569	1574	1586	rBV2	16012	36697	1.70%	0.135%
34	12.607	1589	1593	1600	rVB2	8247	13362	0.62%	0.049%
35	12.966	1651	1654	1661	rVB5	6170	9817	0.46%	0.036%
36	13.172	1684	1689	1695	rBV5	5862	8768	0.41%	0.032%

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Smoothing : OFF

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Min Area: 0.1 % of largest Peak

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Max Peaks: 100

Stop Thrs : 0

Peak Location: TOP

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

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Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM013122.M
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37	13.395	1720	1727	1734	rVB2	26423	39930	1.85%	0.147%
38	13.466	1734	1739	1744	rBV4	6985	12398	0.57%	0.046%
39	13.883	1806	1810	1815	rBV4	8200	14095	0.65%	0.052%
40	13.966	1815	1824	1831	rBV	837333	1116819	51.78%	4.113%

41	14.207	1861	1865	1867	rBV2	19939	29366	1.36%	0.108%
42	14.248	1867	1872	1884	rVB	821675	1249560	57.93%	4.602%
43	14.472	1904	1910	1917	rBV2	29443	67204	3.12%	0.248%
44	14.560	1918	1925	1931	rVV	894890	1392549	64.56%	5.129%
45	14.625	1931	1936	1941	rVB2	14038	23178	1.07%	0.085%

46	14.748	1950	1957	1970	rBV2	168551	343678	15.93%	1.266%
47	14.907	1979	1984	1988	rBV7	6199	10747	0.50%	0.040%
48	14.960	1988	1993	1997	rBV3	15692	28438	1.32%	0.105%
49	15.078	2009	2013	2019	rVB7	5406	9590	0.44%	0.035%
50	15.154	2021	2026	2029	rVV	29123	41373	1.92%	0.152%

51	15.183	2029	2031	2034	rVV5	11158	15249	0.71%	0.056%
52	15.248	2039	2042	2046	rVB3	6993	9210	0.43%	0.034%
53	15.360	2055	2061	2065	rBV3	20157	38054	1.76%	0.140%
54	15.407	2066	2069	2074	rVB3	19230	24574	1.14%	0.091%
55	15.542	2086	2092	2098	rBV	1110881	1586346	73.55%	5.843%

56	15.595	2098	2101	2106	rVV3	20980	39067	1.81%	0.144%
57	15.660	2106	2112	2119	rVV	241421	347514	16.11%	1.280%
58	15.801	2126	2136	2142	rBV6	19695	44245	2.05%	0.163%
59	15.907	2148	2154	2159	rBV7	9624	19921	0.92%	0.073%
60	16.030	2171	2175	2178	rBV6	5729	8302	0.38%	0.031%

61	16.224	2205	2208	2212	rBV6	9144	13106	0.61%	0.048%
62	16.266	2212	2215	2218	rBV2	26321	36229	1.68%	0.133%
63	16.436	2240	2244	2249	rVB7	10162	15131	0.70%	0.056%
64	16.548	2258	2263	2267	rBV2	27008	37916	1.76%	0.140%
65	16.601	2267	2272	2277	rVV	51368	74359	3.45%	0.274%

66	16.666	2277	2283	2286	rBV5	12767	20248	0.94%	0.075%
67	16.836	2309	2312	2315	rBV5	7483	10243	0.47%	0.038%
68	16.954	2327	2332	2340	rBV	24083	51523	2.39%	0.190%
69	17.054	2346	2349	2355	rBV2	31965	45334	2.10%	0.167%
70	17.113	2355	2359	2364	rVB4	32523	49945	2.32%	0.184%

71	17.213	2373	2376	2380	rVB3	18417	23725	1.10%	0.087%
72	17.295	2384	2390	2394	rBV	1002286	1466110	67.97%	5.400%
73	17.336	2394	2397	2401	rVV	155360	201193	9.33%	0.741%
74	17.389	2401	2406	2415	rVV2	1085567	1617744	75.00%	5.958%
75	17.454	2415	2417	2421	rVB2	29883	31180	1.45%	0.115%

76	17.795	2471	2475	2481	rBV	43205	56705	2.63%	0.209%
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 Sampling : 1 Min Area: 0.1 % of largest Peak
 Start Thrs: 0.2 Max Peaks: 100
 Stop Thrs : 0 Peak Location: TOP

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77	17.871	2486	2488	2495	rVB6	9405	11262	0.52%	0.041%
78	17.966	2500	2504	2509	rVV	91342	106442	4.93%	0.392%
79	18.077	2520	2523	2525	rBV3	13227	16654	0.77%	0.061%
80	18.189	2535	2542	2544	rBV2	63332	121688	5.64%	0.448%
81	18.301	2558	2561	2564	rBV2	21984	25349	1.18%	0.093%
82	18.360	2567	2571	2575	rBV4	30997	56665	2.63%	0.209%
83	18.489	2589	2593	2597	rBV	81378	104991	4.87%	0.387%
84	18.666	2621	2623	2627	rBV3	19266	16716	0.77%	0.062%
85	18.954	2668	2672	2677	rBV3	38312	63419	2.94%	0.234%
86	19.118	2696	2700	2706	rBV	114708	230594	10.69%	0.849%
87	19.342	2734	2738	2743	rVB	238443	312193	14.47%	1.150%
88	19.407	2745	2749	2754	rVB	337035	398571	18.48%	1.468%
89	19.677	2790	2795	2806	rBV2	1360662	2156964	100.00%	7.944%
90	19.889	2828	2831	2837	rVB5	48678	65337	3.03%	0.241%
91	20.083	2862	2864	2869	rBV2	47099	51340	2.38%	0.189%
92	20.224	2885	2888	2892	rVB	159573	164435	7.62%	0.606%
93	20.407	2915	2919	2925	rBV5	74012	165098	7.65%	0.608%
94	20.718	2969	2972	2974	rBV	92780	99318	4.60%	0.366%
95	21.177	3047	3050	3056	rVB2	157748	178278	8.27%	0.657%
96	21.324	3072	3075	3081	rBV2	144222	251024	11.64%	0.925%
97	21.459	3093	3098	3102	rBV	1063918	1417985	65.74%	5.222%
98	22.654	3297	3301	3310	rVB	708419	995903	46.17%	3.668%
99	23.695	3472	3478	3483	rBV2	770014	1396894	64.76%	5.145%
100	23.848	3498	3504	3511	rVB2	689853	1334024	61.85%	4.913%

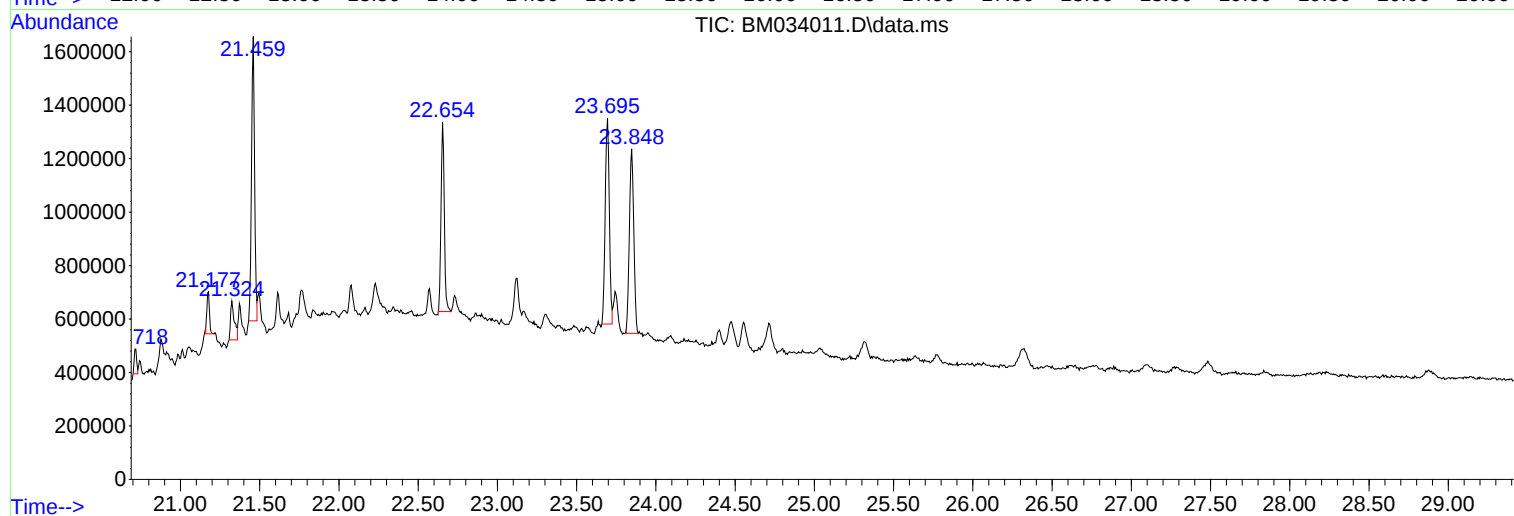
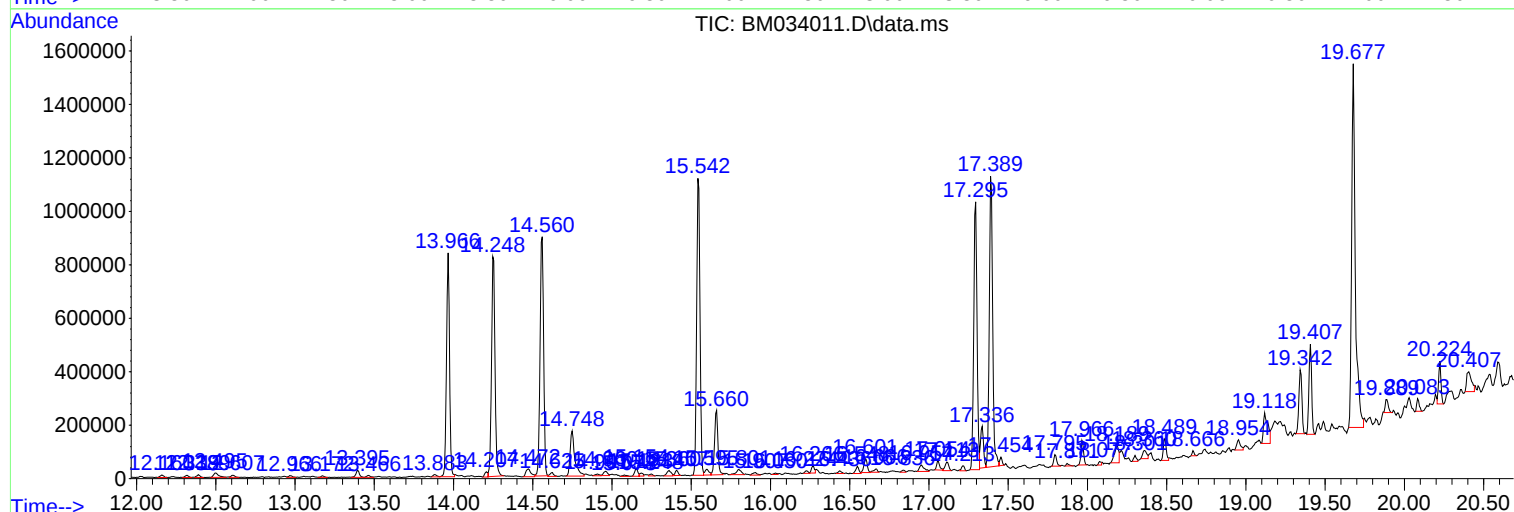
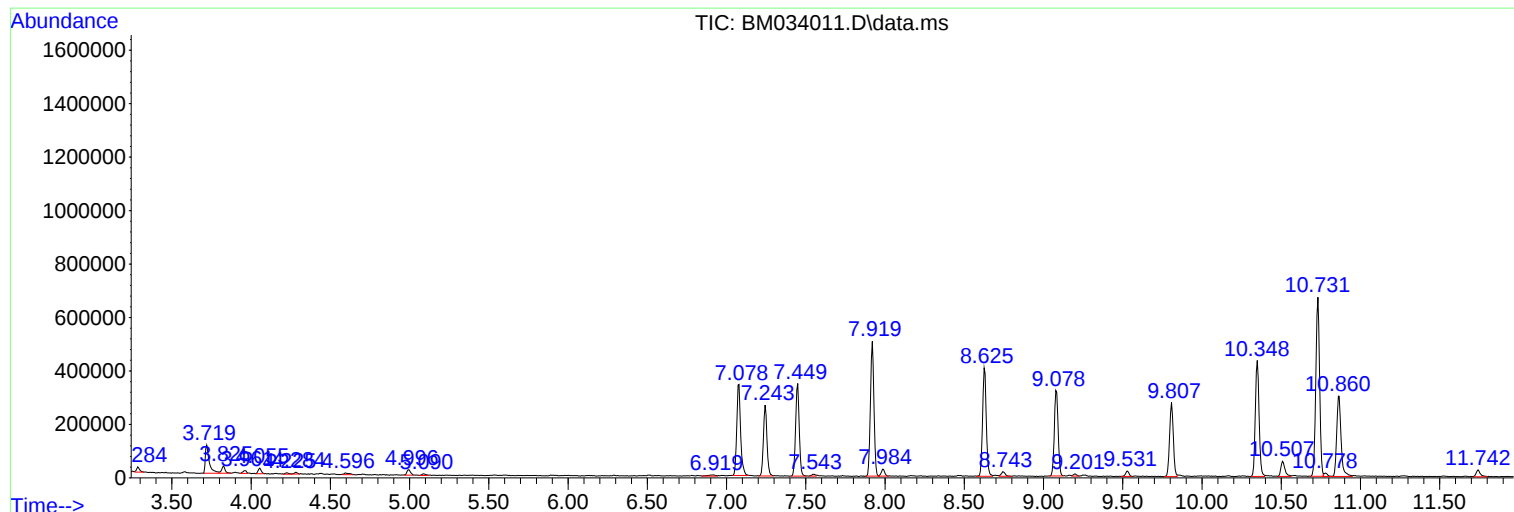
Sum of corrected areas: 27151627

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Instrument :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM013122.M
Quant Title : SVOA CALIBRATION

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



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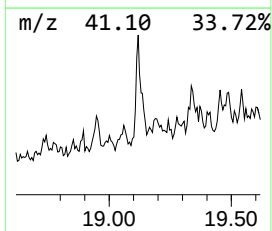
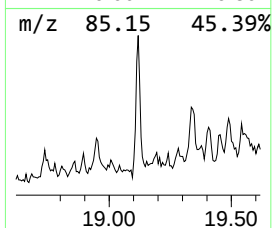
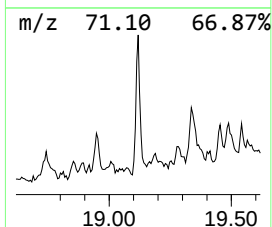
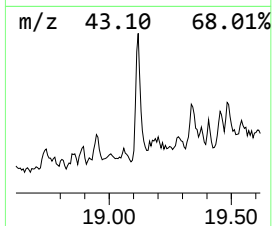
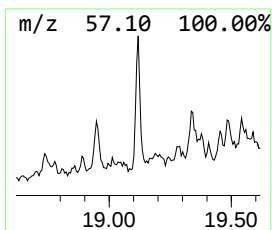
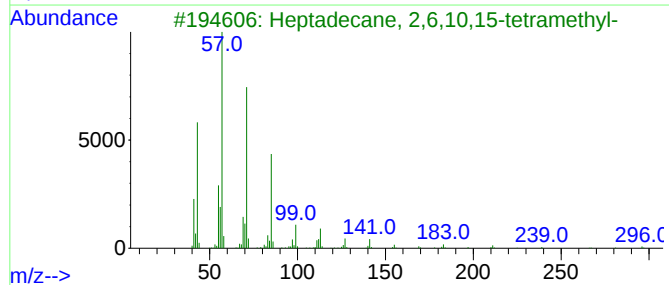
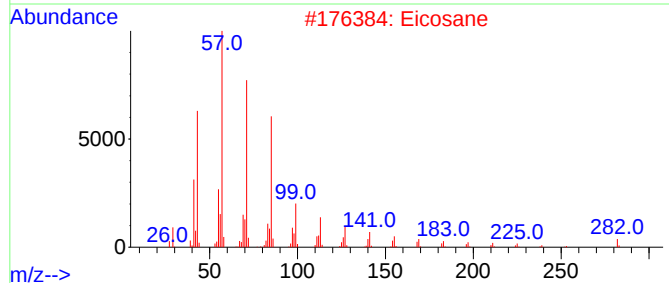
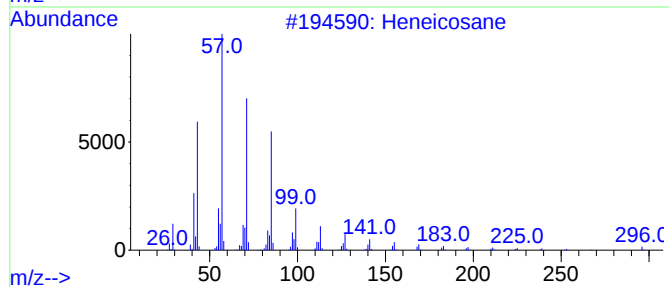
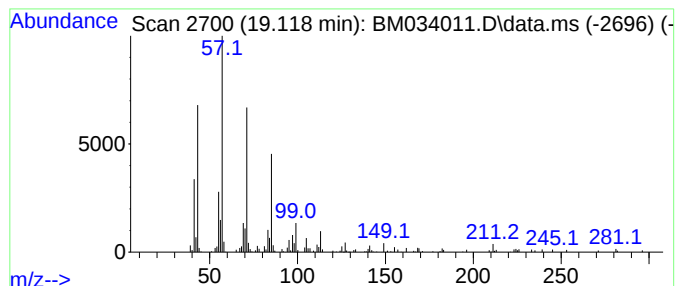
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM013122.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 (DEL) Alkane: Straight-Chai... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.118	3.15 ng/ul	230594	Phenanthrene-d10	17.295

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heneicosane	296	C21H44	000629-94-7	96
2			Eicosane	282	C20H42	000112-95-8	95
3			Heptadecane, 2,6,10,15-tetramethyl-	296	C21H44	054833-48-6	93
4			Heneicosane	296	C21H44	000629-94-7	91
5			Hexacosane	366	C26H54	000630-01-3	87



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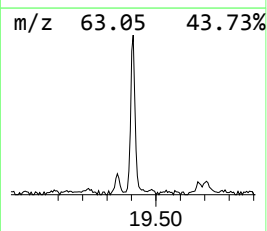
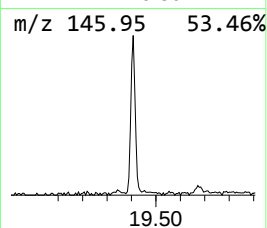
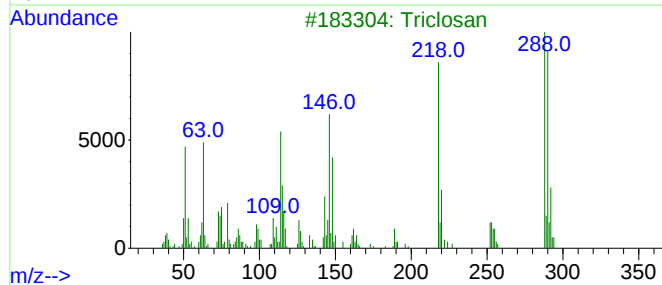
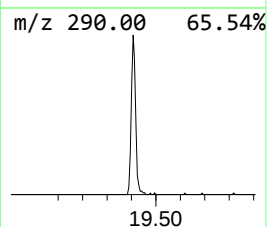
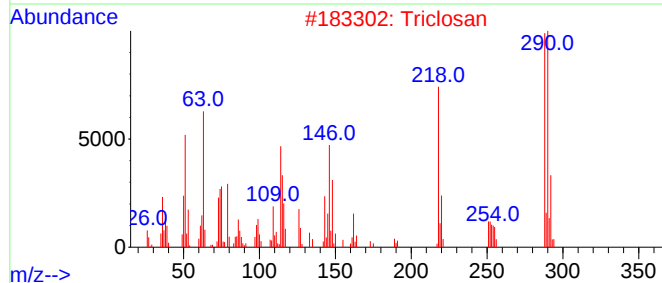
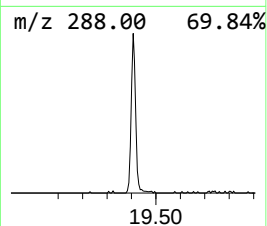
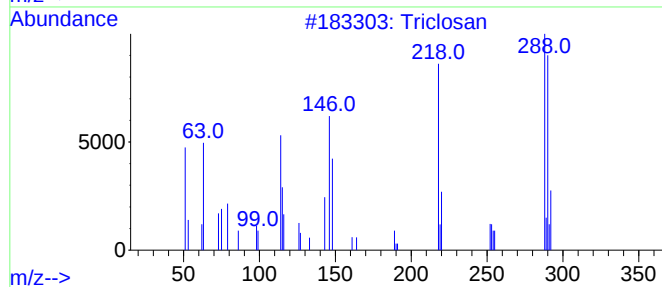
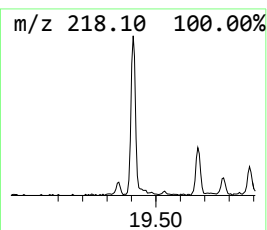
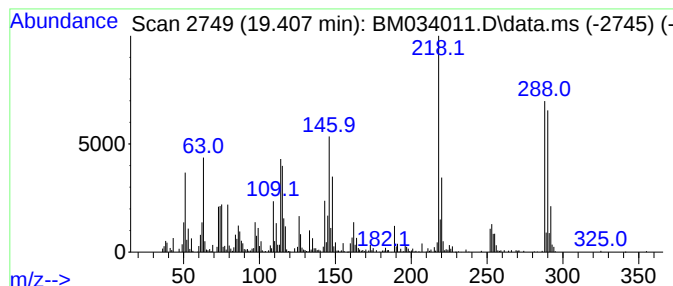
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Quant Title : SVOA CALIBRATION

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

Peak Number 2 Triclosan Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.407	5.62 ng/ul	398571	Chrysene-d12	21.459

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Triclosan	288	C12H7Cl3O2	003380-34-5	99
2			Triclosan	288	C12H7Cl3O2	003380-34-5	99
3			Triclosan	288	C12H7Cl3O2	003380-34-5	91
4			4,4'-Trichlorodiphenylsulfide	288	C12H7Cl3S	1000283-58-7	80
5			Dibenzo[b,e][1,4]dioxin, 1-chloro-	218	C12H7ClO2	039227-53-7	49



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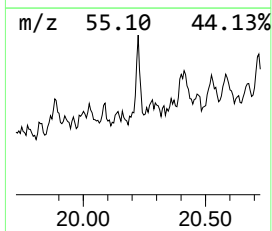
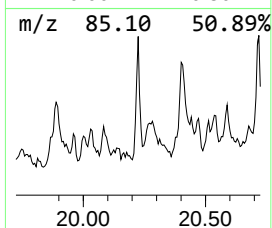
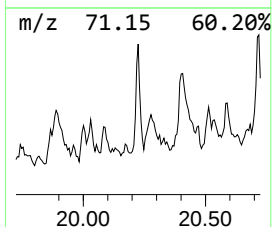
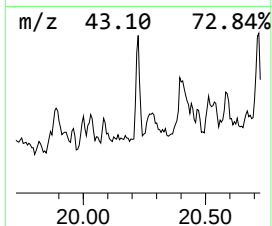
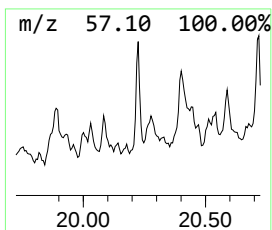
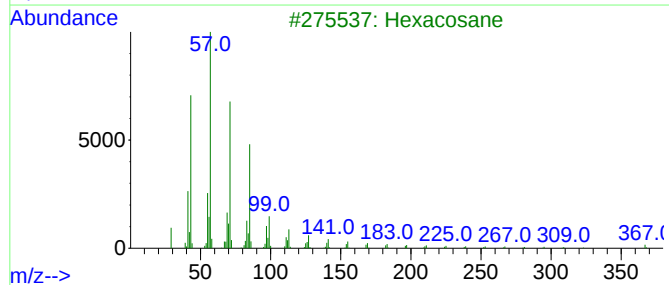
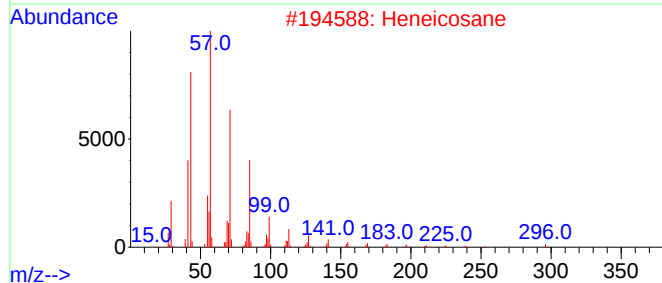
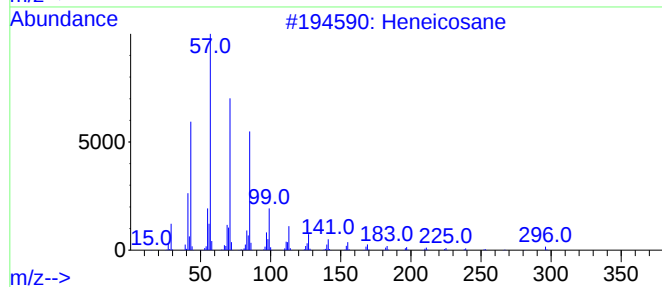
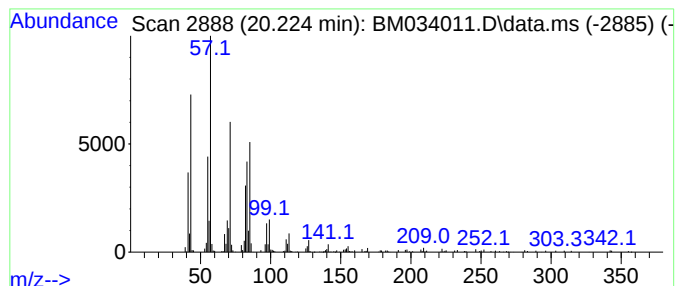
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM013122.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 (DEL) Alkane: Straight-Chai... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.224	2.32 ng/ul	164435	Chrysene-d12	21.459

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heneicosane	296	C21H44	000629-94-7	92
2			Heneicosane	296	C21H44	000629-94-7	83
3			Hexacosane	366	C26H54	000630-01-3	70
4			Dodecane, 1-iodo-	296	C12H25I	004292-19-7	70
5			Heptadecane, 2,6,10,15-tetramethyl-	296	C21H44	054833-48-6	66



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM020322\
Data File : BM034011.D
Acq On : 03 Feb 2022 20:34
Operator : CG/JU
Sample : N1365-02
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
BGLY8

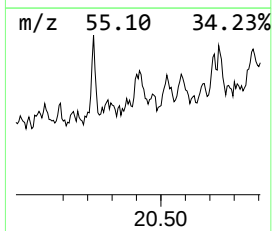
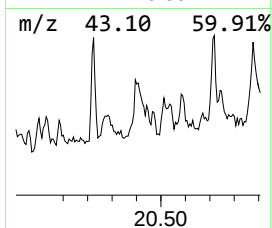
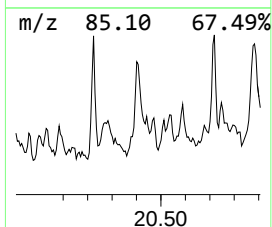
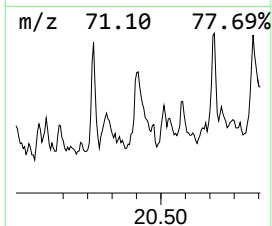
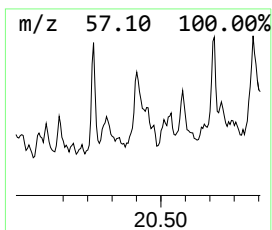
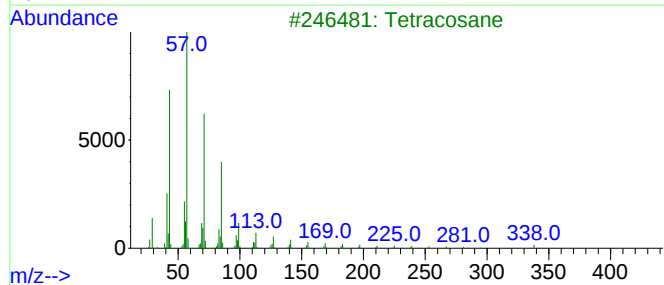
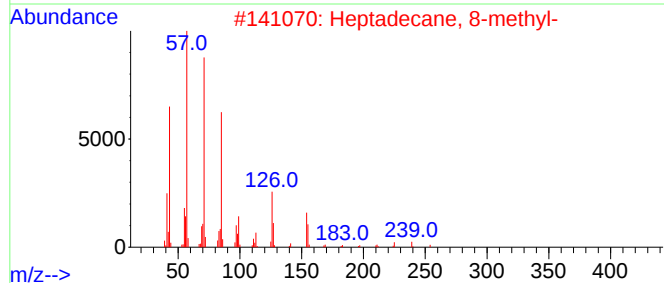
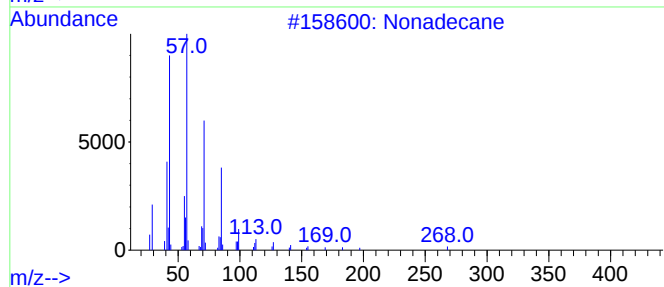
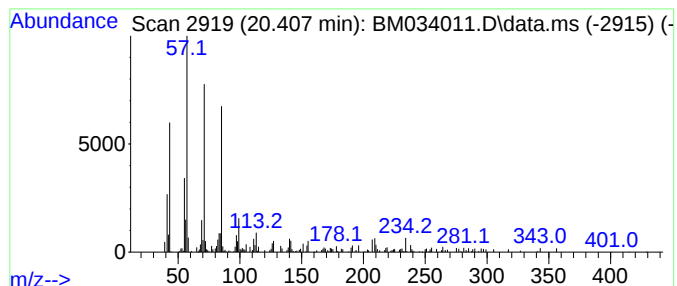
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 (DEL) Alkane: Straight-Chai... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.407	2.33 ng/ul	165098	Chrysene-d12	21.459

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Nonadecane	268	C19H40	000629-92-5	90
2			Heptadecane, 8-methyl-	254	C18H38	013287-23-5	87
3			Tetracosane	338	C24H50	000646-31-1	86
4			Dodecane, 1-iodo-	296	C12H25I	004292-19-7	86
5			Octadecane, 1-iodo-	380	C18H37I	000629-93-6	86



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM020322\
Data File : BM034011.D
Acq On : 03 Feb 2022 20:34
Operator : CG/JU
Sample : N1365-02
Misc :
ALS Vial : 19 Sample Multiplier: 1

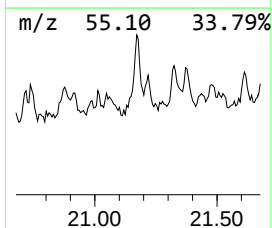
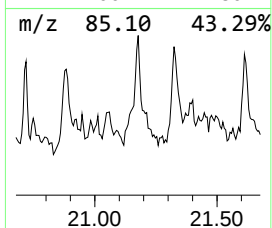
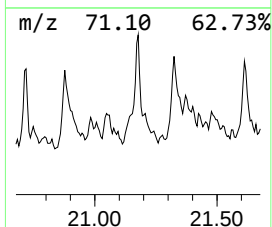
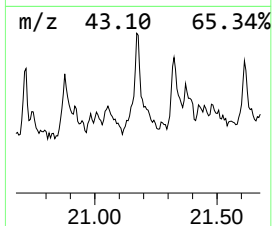
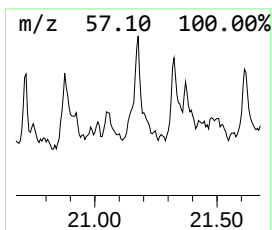
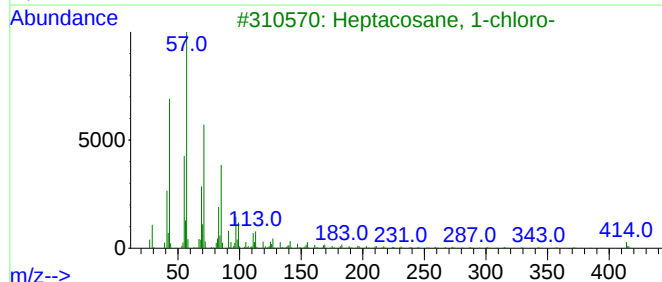
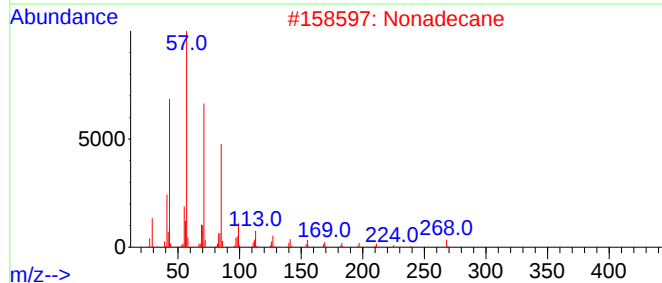
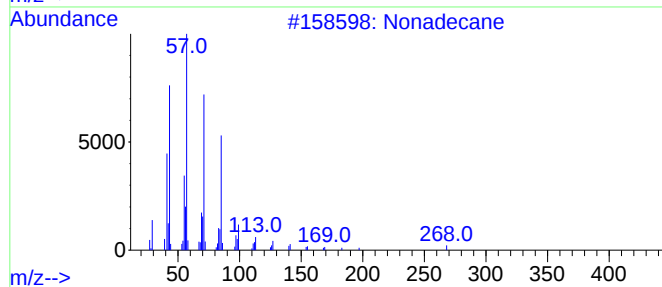
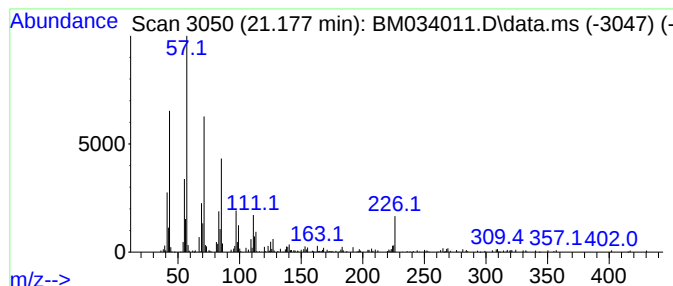
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM013122.M
Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD		R.T.	
21.177	2.51 ng/ul	178278	Chrysene-d12		21.459	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Nonadecane		268	C19H40	000629-92-5	96
2	Nonadecane		268	C19H40	000629-92-5	94
3	Heptacosane, 1-chloro-		414	C27H55Cl	062016-79-9	87
4	Tritetracontane		605	C43H88	007098-21-7	86
5	Tetratetracontane		619	C44H90	007098-22-8	83



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM020322\
Data File : BM034011.D
Acq On : 03 Feb 2022 20:34
Operator : CG/JU
Sample : N1365-02
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
BGLY8

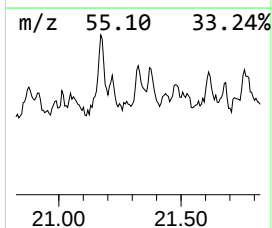
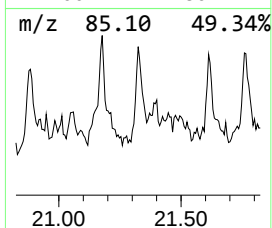
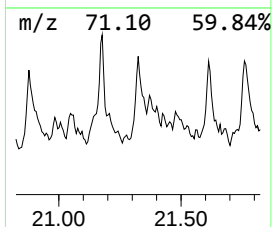
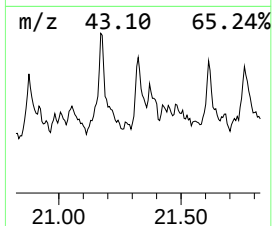
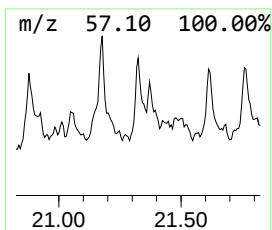
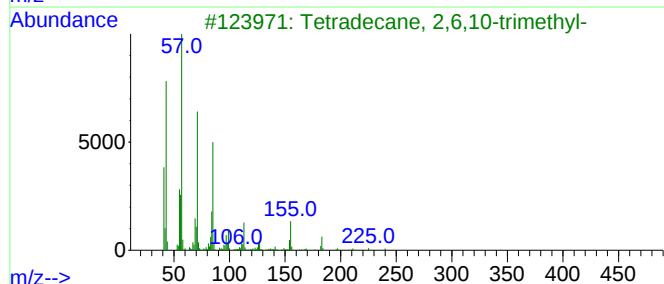
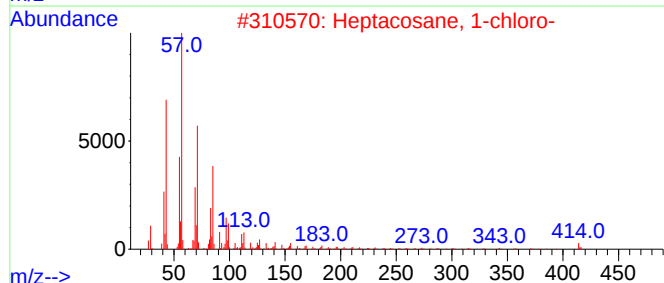
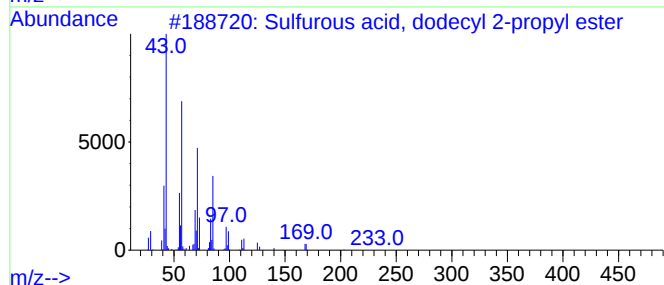
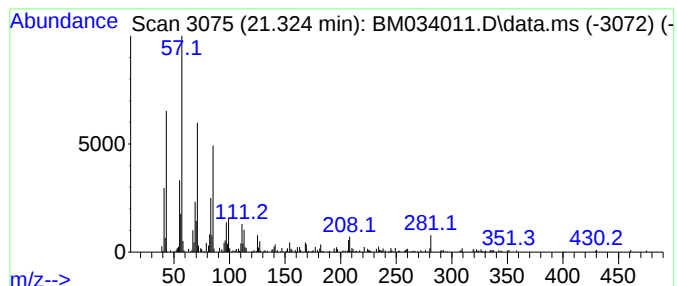
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 Sulfurous acid, dodecyl 2-p... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.324	3.54 ng/ul	251024	Chrysene-d12	21.459

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Sulfurous acid, dodecyl 2-propyl...	292	C15H32O3S	1000309-12-3	87
2			Heptacosane, 1-chloro-	414	C27H55Cl	062016-79-9	86
3			Tetradecane, 2,6,10-trimethyl-	240	C17H36	014905-56-7	83
4			Carbonic acid, eicosyl vinyl ester	368	C23H44O3	1000382-54-3	83
5			Sulfurous acid, pentadecyl 2-pro...	334	C18H38O3S	1000309-12-6	83



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM020322\
Data File : BM034011.D
Acq On : 03 Feb 2022 20:34
Operator : CG/JU
Sample : N1365-02
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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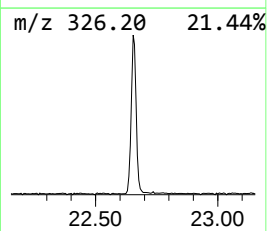
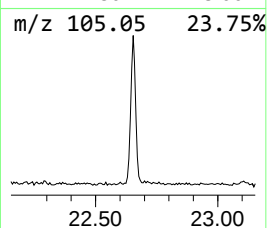
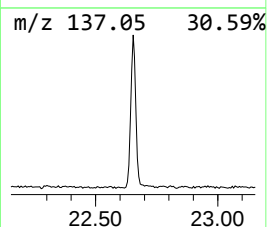
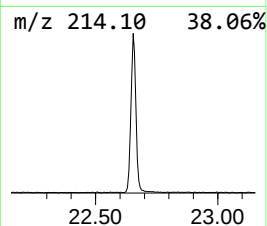
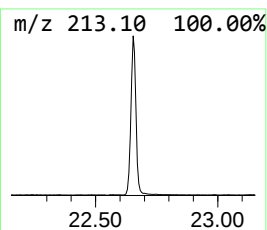
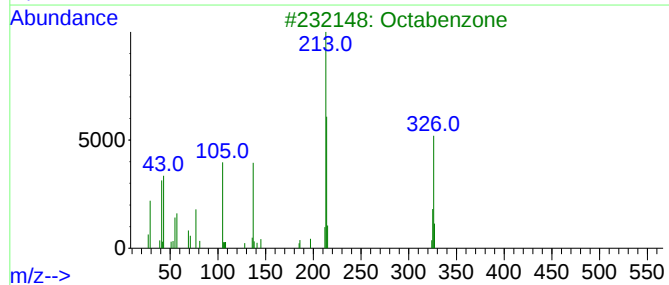
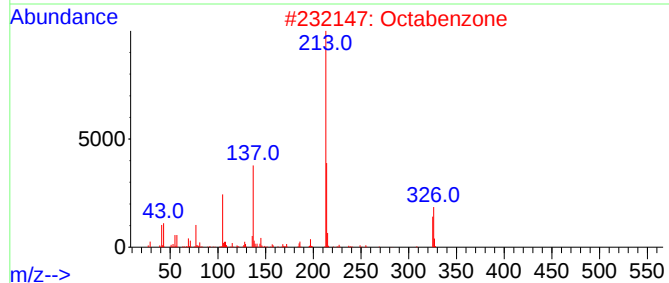
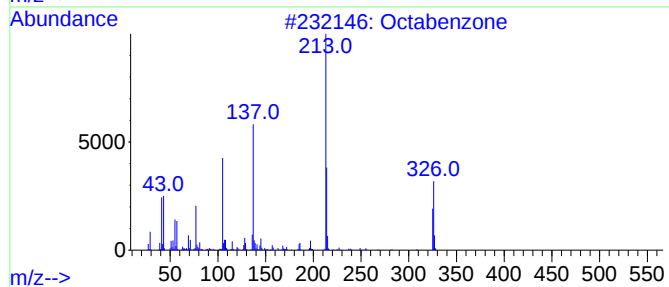
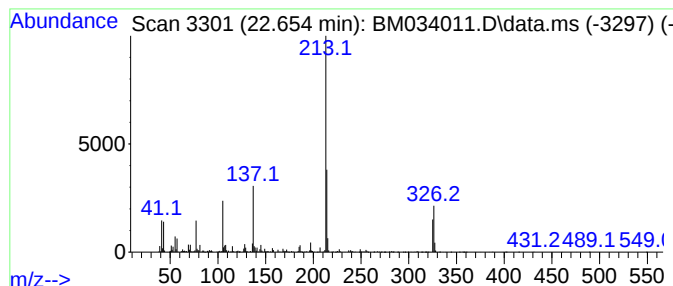
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 Octabenzene Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.654	14.05 ng/ul	995903	Chrysene-d12	21.459

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octabenzene	326	C21H26O3	001843-05-6	98
2			Octabenzene	326	C21H26O3	001843-05-6	94
3			Octabenzene	326	C21H26O3	001843-05-6	93
4			Octabenzene	326	C21H26O3	001843-05-6	52
5			4-Benzoyl-1,3-phenylene diacetate	298	C17H14O5	1000385-78-2	46



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM020322\
Data File : BM034011.D
Acq On : 03 Feb 2022 20:34
Operator : CG/JU
Sample : N1365-02
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
BGLY8

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM013122.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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Triclosan	19.407	5.6	ng/u1	398571	5	21.459	1417990	20.0
(DEL) Alkane: S...	20.224	2.3	ng/u1	164435	5	21.459	1417990	20.0
(DEL) Alkane: S...	20.407	2.3	ng/u1	165098	5	21.459	1417990	20.0
(DEL) Alkane: S...	21.177	2.5	ng/u1	178278	5	21.459	1417990	20.0
Sulfurous acid,...	21.324	3.5	ng/u1	251024	5	21.459	1417990	20.0
Octabenzene	22.654	14.1	ng/u1	995903	5	21.459	1417990	20.0