

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120921\
 Data File : BG051465.D
 Acq On : 10 Dec 2021 21:03
 Operator : CG/JU
 Sample : M4985-12
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 EW5Q6

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_G\Methods\SFAM-EPA-BG120821.M
 Title : SVOA CALIBRATION

Signal : TIC: BG051465.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.530	4	7	13	rVB2	6311	9762	1.15%	0.115%
2	4.000	83	87	91	rBV3	10476	20474	2.40%	0.241%
3	4.252	127	130	131	rBV3	1668	1665	0.20%	0.020%
4	4.299	133	138	148	rVB2	15338	29590	3.47%	0.348%
5	4.487	169	170	173	rBV2	973	982	0.12%	0.012%
6	4.640	194	196	198	rVB3	1078	894	0.10%	0.011%
7	4.687	199	204	212	rBV5	2996	6592	0.77%	0.078%
8	4.840	225	230	233	rBV4	4624	8873	1.04%	0.104%
9	4.869	233	235	243	rVB4	6121	9797	1.15%	0.115%
10	4.928	243	245	248	rBV	1013	911	0.11%	0.011%
11	5.651	364	368	375	rVB	11638	18040	2.12%	0.212%
12	5.786	384	391	404	rBV	36078	87324	10.25%	1.028%
13	6.115	444	447	449	rVB2	846	1030	0.12%	0.012%
14	6.162	449	455	463	rBV	16649	29641	3.48%	0.349%
15	6.291	474	477	480	rBV2	649	918	0.11%	0.011%
16	6.485	507	510	513	rVB3	852	879	0.10%	0.010%
17	6.561	518	523	525	rBV2	1013	1753	0.21%	0.021%
18	7.384	656	663	675	rBV2	13911	34055	4.00%	0.401%
19	7.501	678	683	695	rVV	81618	150877	17.70%	1.776%
20	7.584	695	697	705	rVB5	1505	2918	0.34%	0.034%
21	7.725	715	721	738	rVB	70887	145020	17.01%	1.707%
22	8.183	792	799	810	rVV	89035	157199	18.44%	1.850%
23	8.923	919	925	938	rBV2	48090	111339	13.06%	1.310%
24	9.047	945	946	952	rVB2	1484	2012	0.24%	0.024%
25	9.293	983	988	995	rBV3	3226	6770	0.79%	0.080%
26	9.370	995	1001	1015	rVV	102998	197312	23.15%	2.322%
27	10.098	1117	1125	1137	rBV	77606	159790	18.75%	1.881%
28	10.333	1163	1165	1169	rVB2	828	934	0.11%	0.011%
29	10.656	1215	1220	1238	rBV	73254	192517	22.59%	2.266%
30	11.009	1273	1280	1291	rVB	128549	236740	27.78%	2.786%
31	11.174	1301	1308	1329	rBV	55913	160777	18.86%	1.892%
32	11.661	1388	1391	1396	rVB2	551	951	0.11%	0.011%
33	11.867	1422	1426	1428	rBV2	643	954	0.11%	0.011%
34	12.607	1550	1552	1553	rBV2	1139	1053	0.12%	0.012%
35	12.625	1553	1555	1559	rVB3	1528	1618	0.19%	0.019%
36	13.600	1717	1721	1727	rVB4	4849	7316	0.86%	0.086%

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120921\
 Data File : BG051465.D
 Acq On : 10 Dec 2021 21:03
 Operator : CG/JU
 Sample : M4985-12
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 EW5Q6

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_G\Methods\SFAM-EPA-BG120821.M
 Title : SVOA CALIBRATION

37	14.211	1819	1825	1837	rBV	249585	407206	47.78%	4.793%
38	14.517	1870	1877	1888	rBV	286755	482944	56.66%	5.684%
39	14.669	1898	1903	1907	rVB	8946	11969	1.40%	0.141%
40	14.816	1921	1928	1937	rBV	217578	346467	40.65%	4.078%
41	15.163	1984	1987	1991	rBV3	3290	6213	0.73%	0.073%
42	15.257	2001	2003	2004	rBV2	5692	4739	0.56%	0.056%
43	15.310	2011	2012	2014	rVB	2002	906	0.11%	0.011%
44	15.351	2017	2019	2023	rVB3	1447	1823	0.21%	0.021%
45	15.498	2042	2044	2046	rVV	766	974	0.11%	0.011%
46	15.568	2052	2056	2060	rVV2	2978	5381	0.63%	0.063%
47	15.615	2060	2064	2069	rVB2	7135	10210	1.20%	0.120%
48	15.809	2090	2097	2107	rBV	384978	639403	75.02%	7.526%
49	15.950	2116	2121	2136	rBV2	103800	214238	25.14%	2.522%
50	16.050	2136	2138	2145	rVV7	4707	8060	0.95%	0.095%
51	16.103	2145	2147	2148	rVV2	1742	1245	0.15%	0.015%
52	16.197	2155	2163	2172	rVB	104525	145805	17.11%	1.716%
53	16.273	2172	2176	2178	rVB2	1372	1521	0.18%	0.018%
54	16.385	2193	2195	2200	rBV2	686	871	0.10%	0.010%
55	16.473	2205	2210	2217	rBV3	4087	8116	0.95%	0.096%
56	16.614	2232	2234	2236	rBV2	1106	1117	0.13%	0.013%
57	16.708	2247	2250	2252	rBV3	626	890	0.10%	0.010%
58	16.779	2259	2262	2263	rBV3	1150	870	0.10%	0.010%
59	16.820	2265	2269	2271	rVV4	1605	1611	0.19%	0.019%
60	16.873	2276	2278	2281	rVB3	1361	1385	0.16%	0.016%
61	16.984	2293	2297	2299	rBV3	1020	1375	0.16%	0.016%
62	17.031	2302	2305	2306	rBV3	1033	1084	0.13%	0.013%
63	17.096	2313	2316	2319	rVB5	2037	2137	0.25%	0.025%
64	17.125	2319	2321	2323	rBV2	1200	1394	0.16%	0.016%
65	17.184	2329	2331	2333	rVB3	1305	1087	0.13%	0.013%
66	17.214	2335	2336	2340	rBV4	1146	1365	0.16%	0.016%
67	17.266	2341	2345	2348	rBV3	11976	14686	1.72%	0.173%
68	17.296	2348	2350	2351	rBV2	2865	2132	0.25%	0.025%
69	17.366	2358	2362	2365	rVB4	2807	4142	0.49%	0.049%
70	17.407	2367	2369	2370	rBV2	1554	1047	0.12%	0.012%
71	17.425	2370	2372	2375	rVV4	1478	1837	0.22%	0.022%
72	17.513	2381	2387	2388	rBV5	2964	4669	0.55%	0.055%
73	17.566	2390	2396	2406	rVV	268700	432182	50.71%	5.087%
74	17.666	2406	2413	2423	rBV2	471142	756244	88.73%	8.901%
75	17.736	2423	2425	2427	rBV3	2354	2328	0.27%	0.027%
76	17.872	2444	2448	2453	rBV7	3370	7494	0.88%	0.088%

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120921\
 Data File : BG051465.D
 Acq On : 10 Dec 2021 21:03
 Operator : CG/JU
 Sample : M4985-12
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 EW5Q6

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_G\Methods\SFAM-EPA-BG120821.M
 Title : SVOA CALIBRATION

77	18.007	2467	2471	2475	rVB	17439	21944	2.57%	0.258%
78	18.265	2508	2515	2517	rBV7	6304	13914	1.63%	0.164%
79	18.447	2542	2546	2550	rBV5	5831	8861	1.04%	0.104%
80	18.547	2562	2563	2567	rVB4	5175	4900	0.57%	0.058%
81	18.606	2571	2573	2576	rVV4	5131	5144	0.60%	0.061%
82	18.641	2576	2579	2583	rVB3	13718	18045	2.12%	0.212%
83	18.694	2583	2588	2592	rBV	24816	31508	3.70%	0.371%
84	19.317	2691	2694	2697	rBV	26052	29521	3.46%	0.347%
85	19.346	2697	2699	2703	rVB4	13548	13842	1.62%	0.163%
86	19.511	2724	2727	2728	rBV3	7114	6656	0.78%	0.078%
87	19.581	2737	2739	2747	rVB4	7658	14717	1.73%	0.173%
88	19.852	2783	2785	2787	rBV3	8931	7564	0.89%	0.089%
89	19.893	2788	2792	2796	rVV	30337	41143	4.83%	0.484%
90	19.946	2796	2801	2811	rVB	570331	852317	100.00%	10.032%
91	20.422	2879	2882	2886	rBV3	11831	16520	1.94%	0.194%
92	20.598	2909	2912	2917	rBV4	9955	21170	2.48%	0.249%
93	20.880	2956	2960	2964	rVB	54684	61974	7.27%	0.729%
94	20.921	2964	2967	2969	rBV3	11456	11727	1.38%	0.138%
95	21.696	3094	3099	3105	rVB	111230	158910	18.64%	1.870%
96	21.867	3122	3128	3142	rVB	218929	432794	50.78%	5.094%
97	23.030	3320	3326	3334	rBV	70636	130070	15.26%	1.531%
98	25.028	3656	3666	3682	rBV2	253230	772128	90.59%	9.088%
99	25.269	3697	3707	3721	rVB2	124054	413649	48.53%	4.869%
100	32.178	4875	4883	4884	rBV	33406	68557	8.04%	0.807%

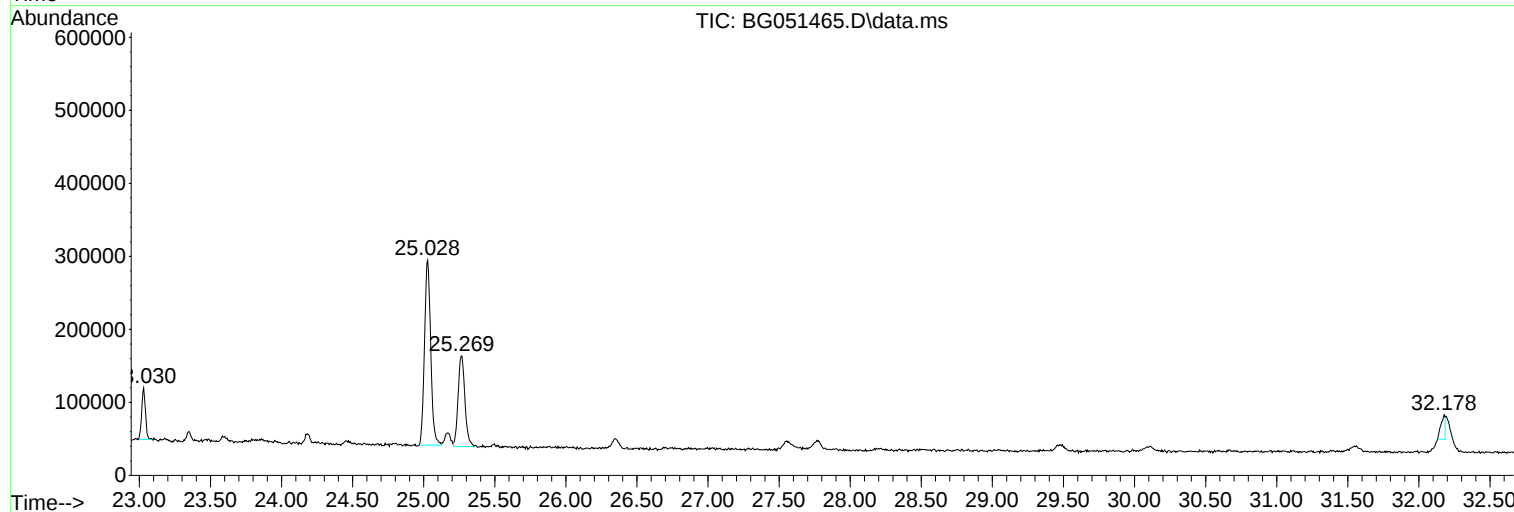
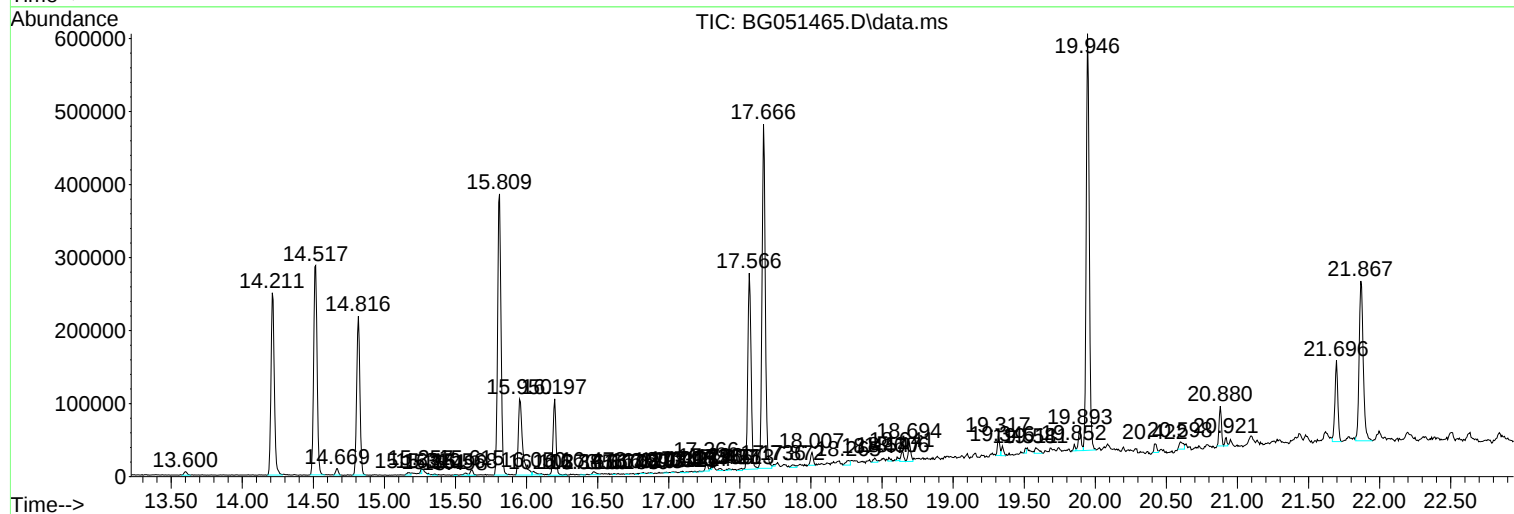
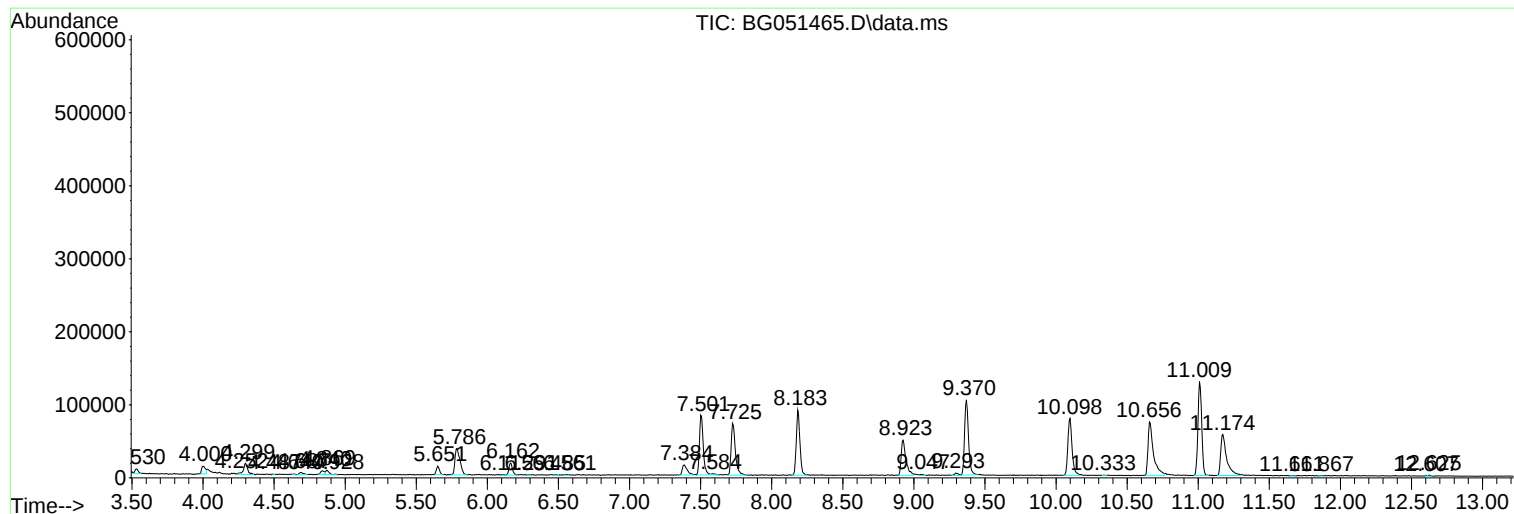
Sum of corrected areas: 8496019

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120921\
Data File : BG051465.D
Acq On : 10 Dec 2021 21:03
Operator : CG/JU
Sample : M4985-12
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
EW5Q6

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG120821.M
Quant Title : SVOA CALIBRATION

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



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120921\
Data File : BG051465.D
Acq On : 10 Dec 2021 21:03
Operator : CG/JU
Sample : M4985-12
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
EW5Q6

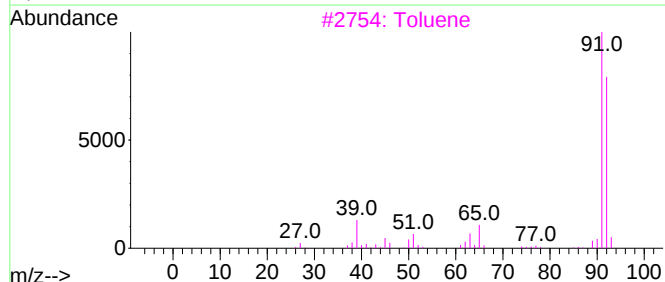
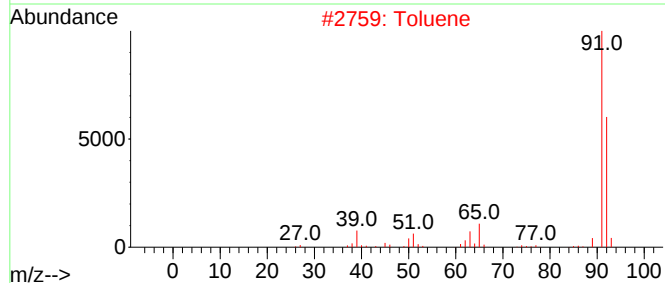
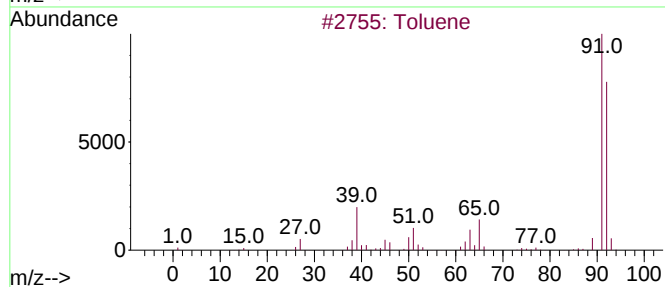
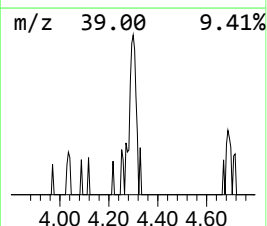
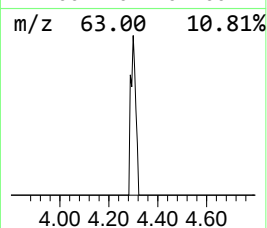
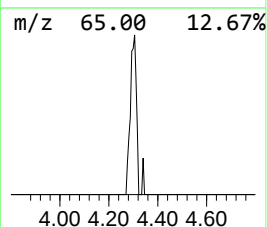
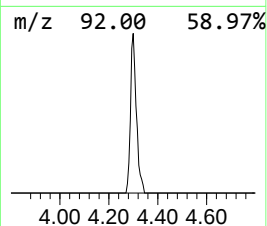
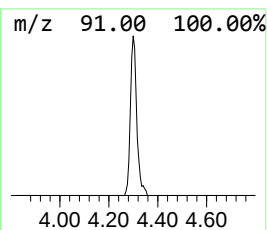
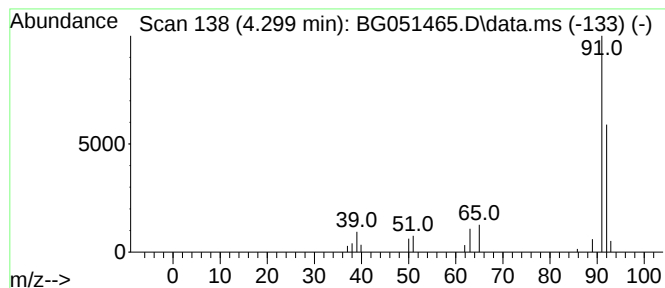
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG120821.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Toluene Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.299	3.76 ng/ul	29590	1,4-Dichlorobenzene-d4	8.183

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Toluene	92	C7H8	000108-88-3	91
2	Toluene	92	C7H8	000108-88-3	90
3	Toluene	92	C7H8	000108-88-3	90
4	Toluene	92	C7H8	000108-88-3	90
5	Toluene	92	C7H8	000108-88-3	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120921\
Data File : BG051465.D
Acq On : 10 Dec 2021 21:03
Operator : CG/JU
Sample : M4985-12
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
EW5Q6

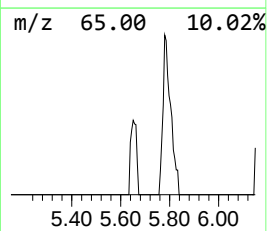
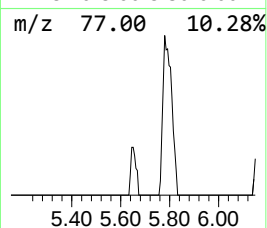
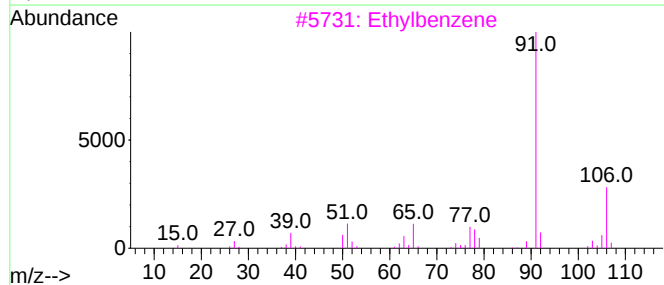
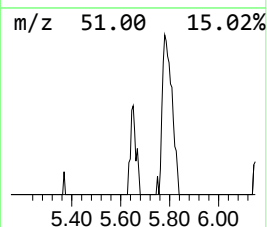
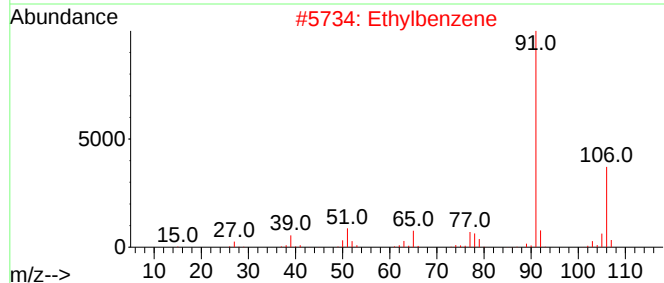
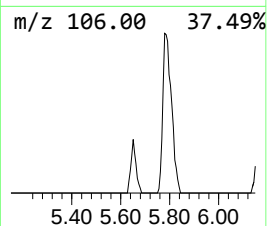
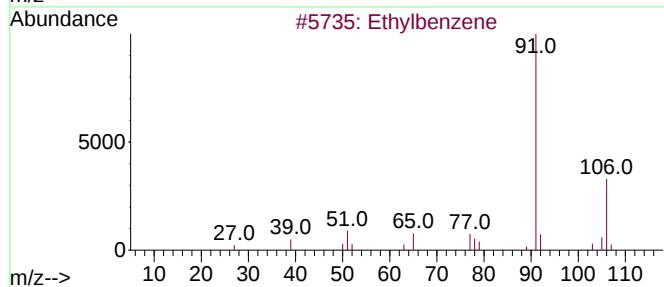
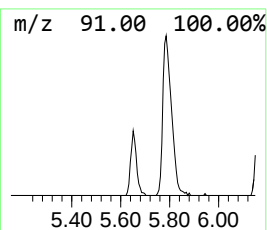
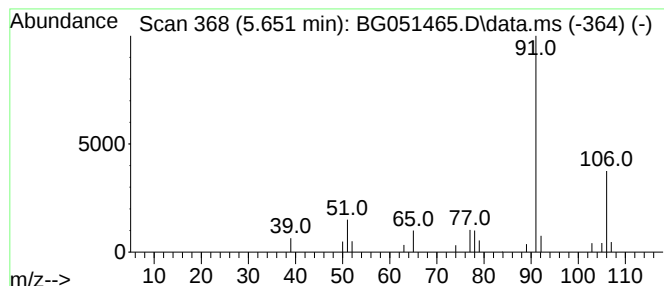
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG120821.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 Ethylbenzene Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.651	2.30 ng/ul	18040	1,4-Dichlorobenzene-d4	8.183

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Ethylbenzene	106	C8H10	000100-41-4	91
2			Ethylbenzene	106	C8H10	000100-41-4	91
3			Ethylbenzene	106	C8H10	000100-41-4	91
4			Ethylbenzene	106	C8H10	000100-41-4	91
5			Benzene, 1,3-dimethyl-	106	C8H10	000108-38-3	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120921\
Data File : BG051465.D
Acq On : 10 Dec 2021 21:03
Operator : CG/JU
Sample : M4985-12
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
EW5Q6

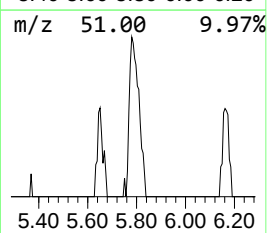
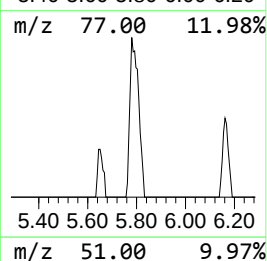
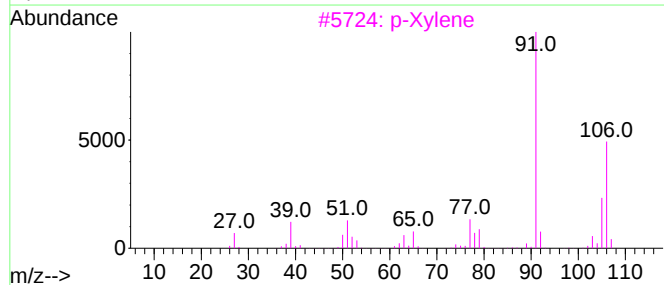
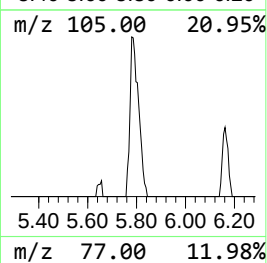
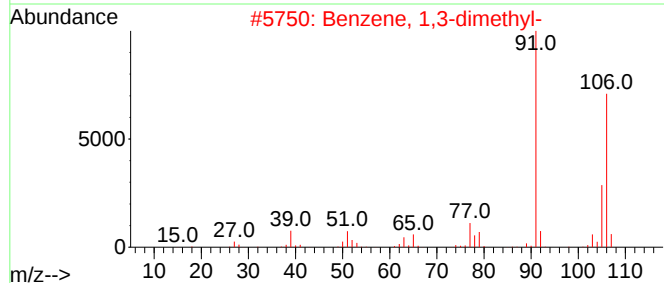
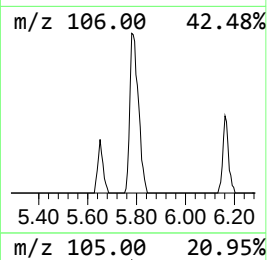
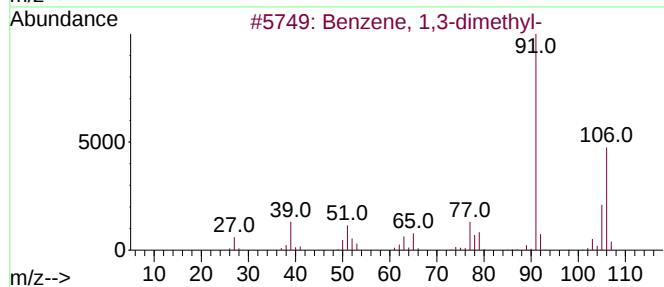
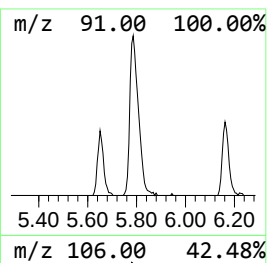
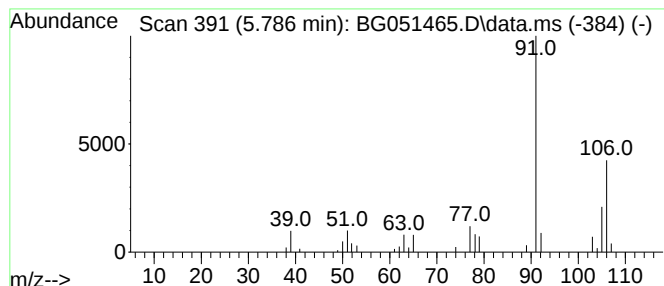
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG120821.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 Benzene, 1,3-dimethyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.786	11.11 ng/ul	87324	1,4-Dichlorobenzene-d4	8.183

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,3-dimethyl-	106	C8H10	000108-38-3	97
2			Benzene, 1,3-dimethyl-	106	C8H10	000108-38-3	95
3			p-Xylene	106	C8H10	000106-42-3	94
4			p-Xylene	106	C8H10	000106-42-3	94
5			Benzene, 1,3-dimethyl-	106	C8H10	000108-38-3	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120921\
Data File : BG051465.D
Acq On : 10 Dec 2021 21:03
Operator : CG/JU
Sample : M4985-12
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
EW5Q6

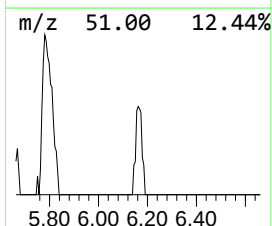
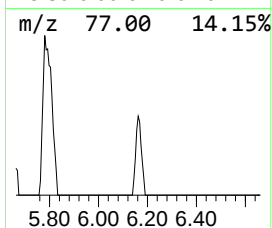
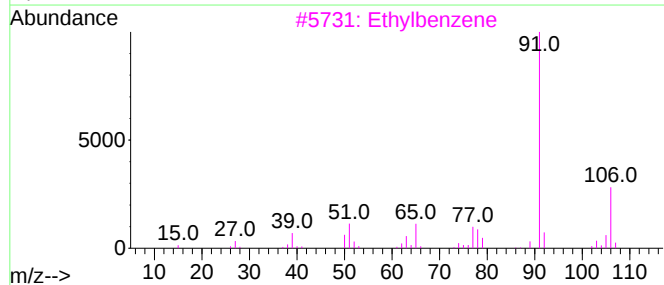
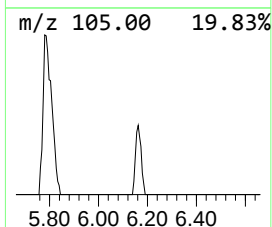
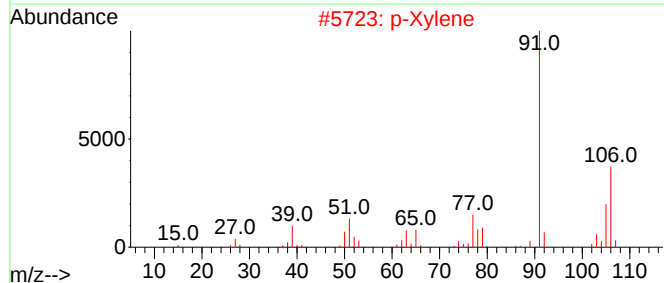
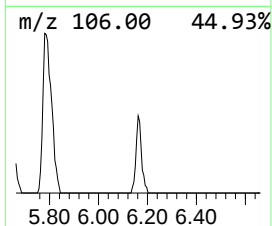
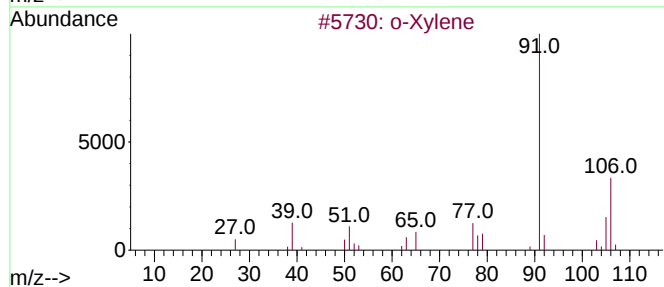
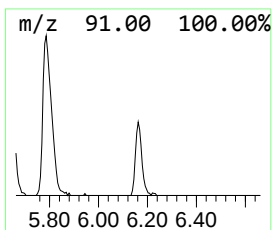
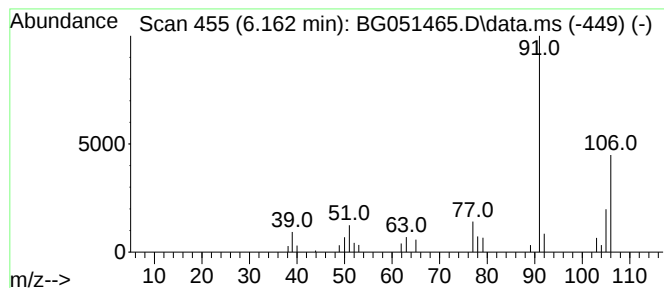
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG120821.M
Quant Title : SVOA CALIBRATION

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

Peak Number 4 o-Xylene Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.162	3.77 ng/ul	29641	1,4-Dichlorobenzene-d4	8.183

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			o-Xylene	106	C8H10	000095-47-6	91
2			p-Xylene	106	C8H10	000106-42-3	90
3			Ethylbenzene	106	C8H10	000100-41-4	72
4			Ethylbenzene	106	C8H10	000100-41-4	72
5			Ethylbenzene	106	C8H10	000100-41-4	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120921\
Data File : BG051465.D
Acq On : 10 Dec 2021 21:03
Operator : CG/JU
Sample : M4985-12
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
EW5Q6

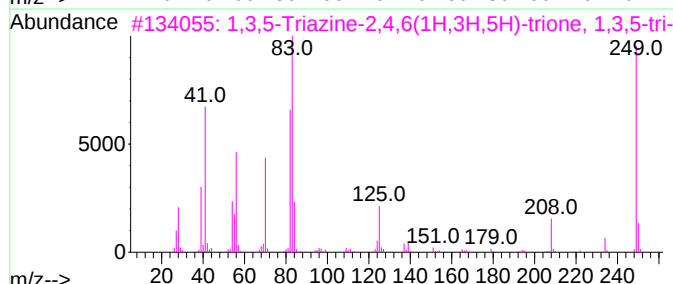
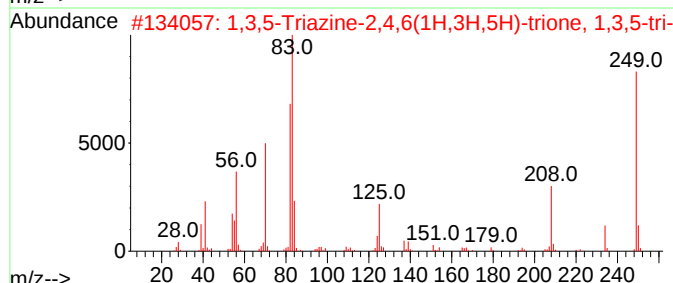
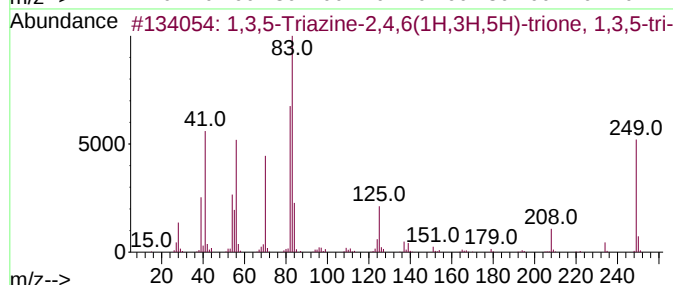
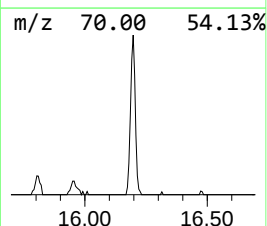
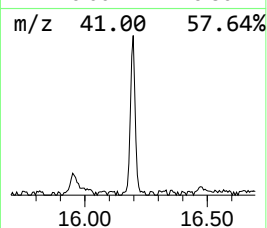
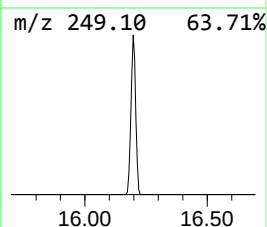
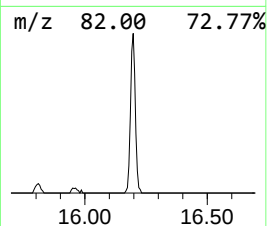
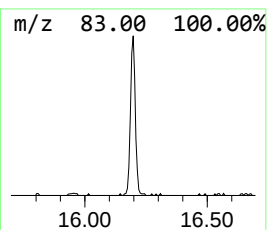
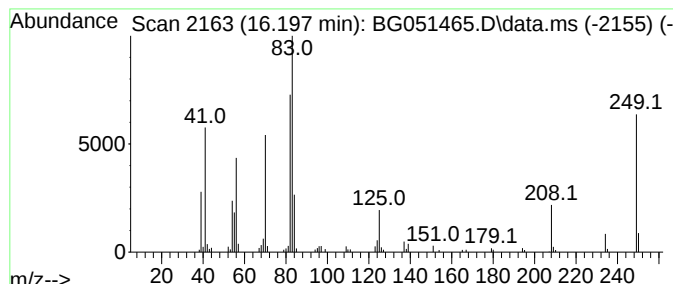
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG120821.M
Quant Title : SVOA CALIBRATION

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

Peak Number 5 1,3,5-Triazine-2,4,6(1H,3H,... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.197	6.75 ng/ul	145805	Phenanthrene-d10	17.566

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1,3,5-Triazine-2,4,6(1H,3H,5H)-t...	249	C12H15N3O3	001025-15-6	98
2			1,3,5-Triazine-2,4,6(1H,3H,5H)-t...	249	C12H15N3O3	001025-15-6	98
3			1,3,5-Triazine-2,4,6(1H,3H,5H)-t...	249	C12H15N3O3	001025-15-6	96
4			1,3,5-Triazine-2,4,6(1H,3H,5H)-t...	249	C12H15N3O3	001025-15-6	95
5			12-Octadecenal	266	C18H34O	056554-91-7	35



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120921\
Data File : BG051465.D
Acq On : 10 Dec 2021 21:03
Operator : CG/JU
Sample : M4985-12
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
EW5Q6

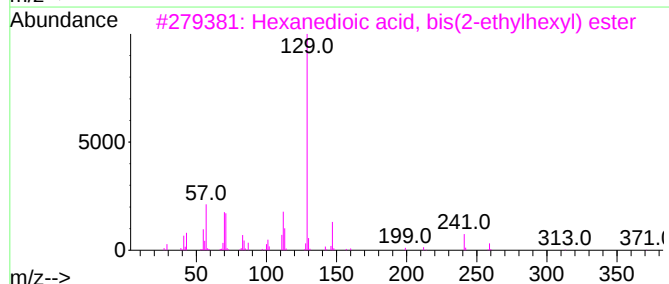
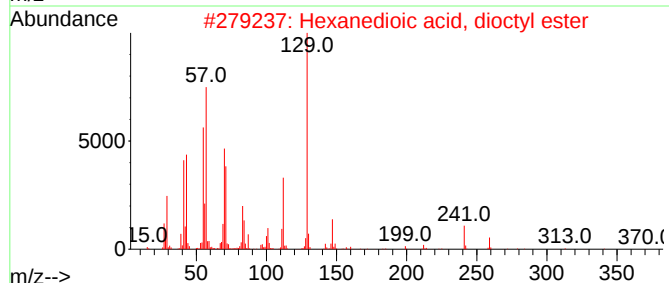
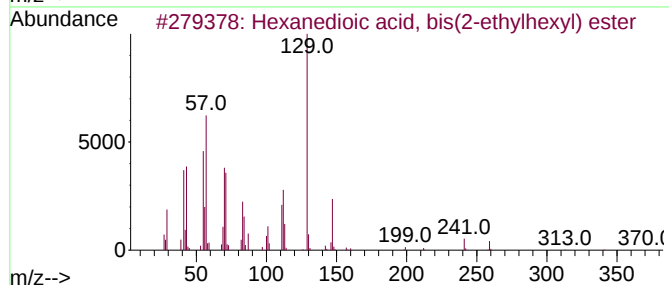
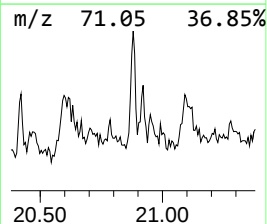
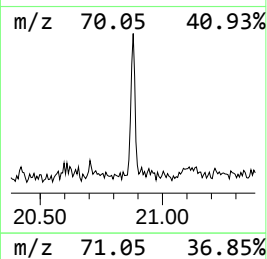
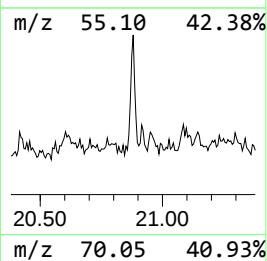
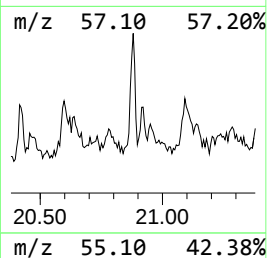
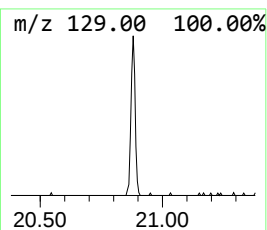
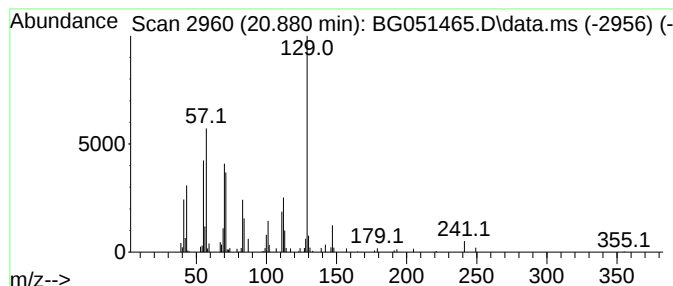
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG120821.M
Quant Title : SVOA CALIBRATION

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

Peak Number 6 Hexanedioic acid, bis(2-eth... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.880	2.86 ng/ul	61974	Chrysene-d12	21.867

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexanedioic acid, bis(2-ethylhex...	370	C22H42O4	000103-23-1	90
2			Hexanedioic acid, dioctyl ester	370	C22H42O4	000123-79-5	87
3			Hexanedioic acid, bis(2-ethylhex...	370	C22H42O4	000103-23-1	78
4			Hexanedioic acid, dioctyl ester	370	C22H42O4	000123-79-5	72
5			Hexanedioic acid, bis(2-ethylhex...	370	C22H42O4	000103-23-1	59



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120921\
Data File : BG051465.D
Acq On : 10 Dec 2021 21:03
Operator : CG/JU
Sample : M4985-12
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
EW5Q6

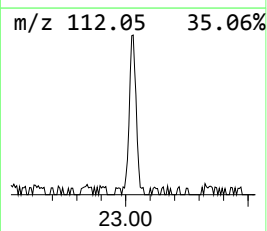
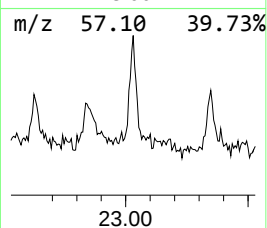
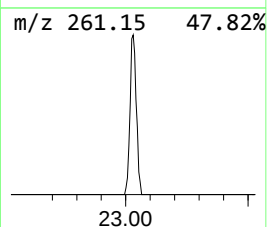
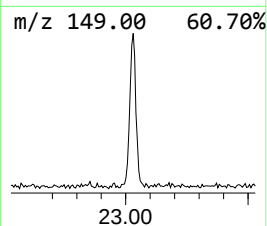
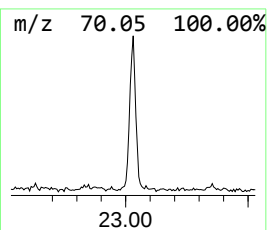
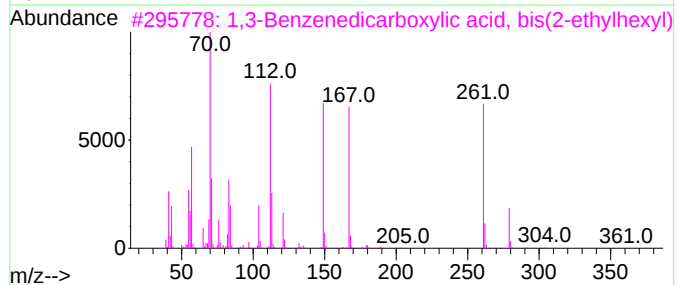
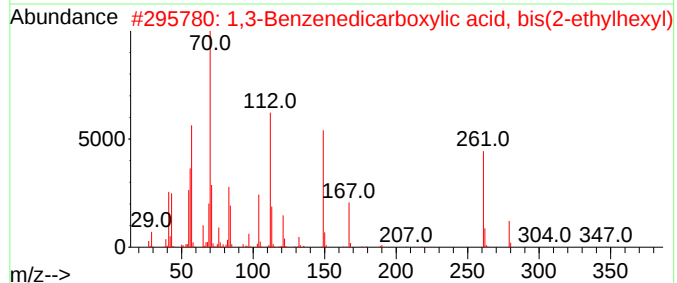
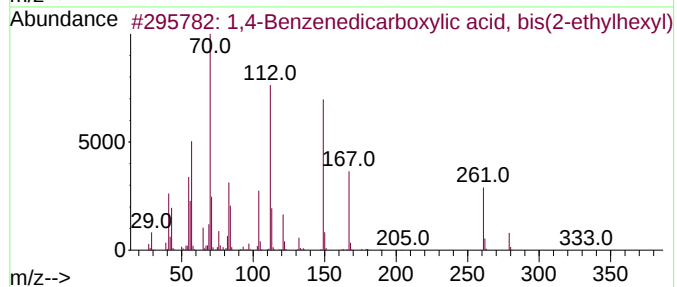
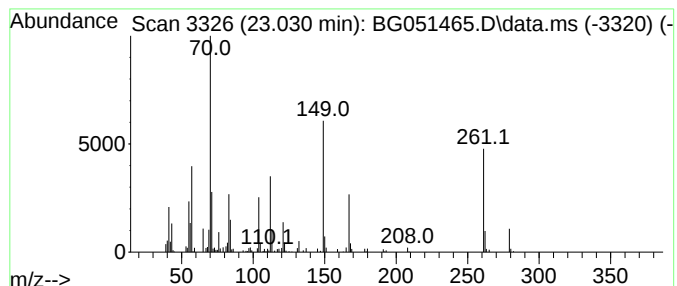
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG120821.M
Quant Title : SVOA CALIBRATION

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

Peak Number 7 1,4-Benzenedicarboxylic aci... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.030	6.01 ng/ul	130070	Chrysene-d12	21.867

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1,4-Benzenedicarboxylic acid, bi...	390	C24H38O4	006422-86-2	59
2			1,3-Benzenedicarboxylic acid, bi...	390	C24H38O4	000137-89-3	49
3			1,3-Benzenedicarboxylic acid, bi...	390	C24H38O4	000137-89-3	47
4			1,4-Benzenedicarboxylic acid, bi...	390	C24H38O4	006422-86-2	38
5			Bis(2-ethylhexyl) phthalate	390	C24H38O4	000117-81-7	22



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120921\
Data File : BG051465.D
Acq On : 10 Dec 2021 21:03
Operator : CG/JU
Sample : M4985-12
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
EW5Q6

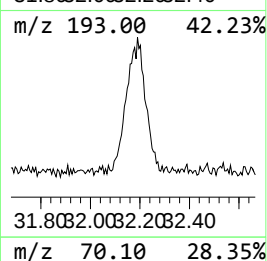
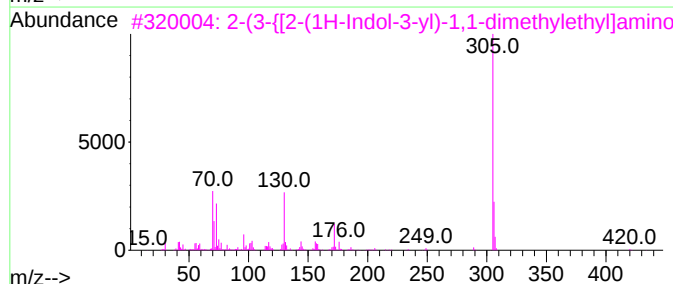
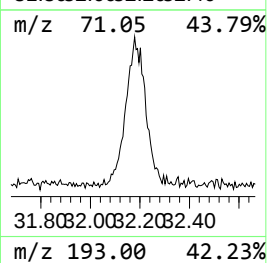
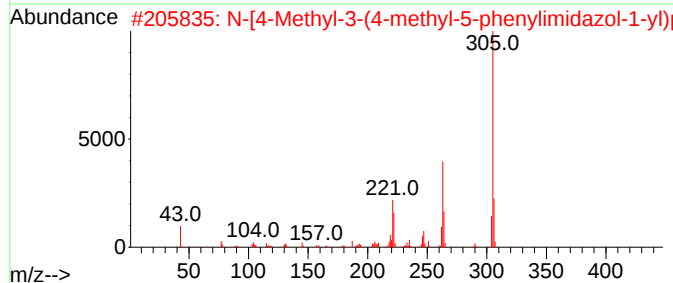
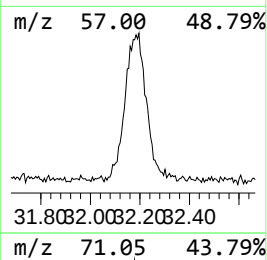
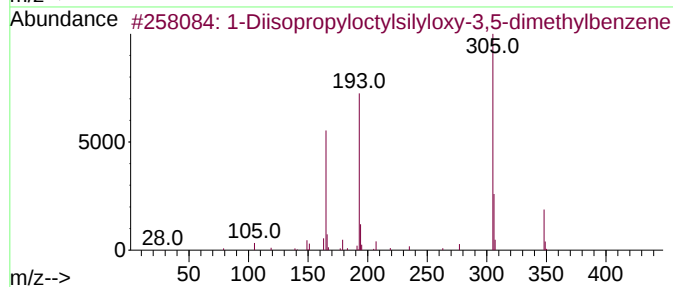
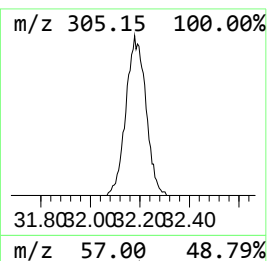
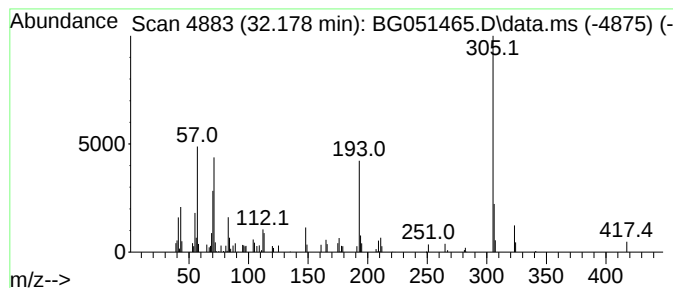
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG120821.M
Quant Title : SVOA CALIBRATION

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

Peak Number 8 1-Diisopropyloctylsilyloxy-... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
32.178	3.31 ng/ul	68557	Perylene-d12	25.263

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Diisopropyloctylsilyloxy-3,5-d...	348	C22H40OSi	1000307-85-1	64		
2	N-[4-Methyl-3-(4-methyl-5-phenyl...	305	C19H19N3O	1000459-63-7	38		
3	2-(3-{[2-(1H-Indol-3-yl)-1,1-dim...	435	C25H33N3O2Si	1000408-41-9	32		
4	(1-pentyl-1H-indol-3-yl)(o-tolyl...	305	C21H23NO	1000439-58-8	25		
5	Bucindolol, 2TMS derivative	507	C28H41N3O2Si2	959296-85-6	25		



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120921\
Data File : BG051465.D
Acq On : 10 Dec 2021 21:03
Operator : CG/JU
Sample : M4985-12
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
EW5Q6

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG120821.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
Toluene	4.299	3.8	ng/ul	29590	1	8.183	157199	20.0
Ethylbenzene	5.651	2.3	ng/ul	18040	1	8.183	157199	20.0
Benzene, 1,3-di...	5.786	11.1	ng/ul	87324	1	8.183	157199	20.0
o-Xylene	6.162	3.8	ng/ul	29641	1	8.183	157199	20.0
1,3,5-Triazine-...	16.197	6.8	ng/ul	145805	4	17.566	432182	20.0
Hexanedioic aci...	20.880	2.9	ng/ul	61974	5	21.867	432794	20.0
1,4-Benzenedica...	23.030	6.0	ng/ul	130070	5	21.867	432794	20.0
1-Diisopropyloc...	32.178	3.3	ng/ul	68557	6	25.263	413649	20.0