

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG102920\
 Data File : BG047118.D
 Acq On : 29 Oct 2020 13:43
 Operator : CG/JU
 Sample : L4499-06
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 EW2B7

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_G\METHODS\SOM-EPA-BG102220MA.M
 Title : SVOA CALIBRATION

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.468	18	22	25	rBV3	5804	8660	0.78%	0.068%
2	3.544	34	35	37	rBV2	3117	2112	0.19%	0.017%
3	3.891	90	94	98	rBV4	8020	11597	1.04%	0.091%
4	3.997	107	112	120	rBV	32601	54773	4.91%	0.431%
5	4.067	120	124	128	rBV4	3845	6071	0.54%	0.048%
6	4.408	180	182	185	rVB4	2391	2075	0.19%	0.016%
7	4.443	185	188	191	rBV4	2765	3863	0.35%	0.030%
8	4.590	210	213	215	rBV2	1999	2282	0.20%	0.018%
9	4.784	242	246	249	rBV5	4228	5569	0.50%	0.044%
10	5.043	289	290	295	rVB3	2966	2701	0.24%	0.021%
11	5.137	303	306	312	rBV6	2195	3486	0.31%	0.027%
12	5.225	317	321	322	rBV	7088	6094	0.55%	0.048%
13	5.366	344	345	350	rBV3	3866	3471	0.31%	0.027%
14	5.524	370	372	374	rBV	1838	2022	0.18%	0.016%
15	5.671	396	397	400	rBV2	2294	2400	0.22%	0.019%
16	6.059	458	463	470	rBV6	2894	6071	0.54%	0.048%
17	6.171	478	482	485	rBV5	1593	2127	0.19%	0.017%
18	7.211	650	659	671	rBV	125118	225905	20.26%	1.776%
19	7.381	679	688	695	rBV	97446	164018	14.71%	1.289%
20	7.534	712	714	716	rBV2	2042	2078	0.19%	0.016%
21	7.587	717	723	730	rVV2	125465	224037	20.09%	1.761%
22	7.640	730	732	735	rVV4	2318	2486	0.22%	0.020%
23	7.875	769	772	775	rBV3	1719	2060	0.18%	0.016%
24	8.057	797	803	811	rVV	224919	381563	34.21%	2.999%
25	8.756	916	922	933	rBV2	129731	238527	21.39%	1.875%
26	9.220	993	1001	1010	rBV	126561	227694	20.42%	1.790%
27	9.514	1049	1051	1057	rVB4	1629	2511	0.23%	0.020%
28	9.943	1117	1124	1133	rBV2	113826	204140	18.30%	1.605%
29	10.483	1209	1216	1226	rBV	163369	300184	26.92%	2.360%
30	10.754	1260	1262	1267	rVB	1726	2196	0.20%	0.017%
31	10.871	1271	1282	1289	rBV	297517	526560	47.21%	4.139%
32	11.000	1296	1304	1313	rBV3	89808	172632	15.48%	1.357%
33	11.723	1424	1427	1429	rBV	1849	2239	0.20%	0.018%
34	11.770	1432	1435	1441	rVV5	1528	2835	0.25%	0.022%

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG102920\
 Data File : BG047118.D
 Acq On : 29 Oct 2020 13:43
 Operator : CG/JU
 Sample : L4499-06
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 EW2B7

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_G\METHODS\SOM-EPA-BG102220MA.M
 Title : SVOA CALIBRATION

35	11.929	1460	1462	1465	rBV2	2165	2324	0.21%	0.018%
36	12.446	1548	1550	1555	rVB3	2947	3747	0.34%	0.029%
37	12.516	1559	1562	1563	rBV2	2245	2550	0.23%	0.020%
38	12.751	1597	1602	1604	rVV4	2305	4039	0.36%	0.032%
39	12.998	1642	1644	1648	rVV	2146	3097	0.28%	0.024%
40	13.568	1738	1741	1747	rBV3	4673	6370	0.57%	0.050%
41	13.644	1749	1754	1756	rBV2	2203	2166	0.19%	0.017%
42	13.709	1762	1765	1768	rBV3	1492	2026	0.18%	0.016%
43	13.862	1789	1791	1793	rBV	2288	2278	0.20%	0.018%
44	14.009	1809	1816	1820	rBV5	2311	5298	0.48%	0.042%
45	14.085	1822	1829	1834	rVV	348280	503933	45.18%	3.961%
46	14.132	1834	1837	1845	rVB	107449	151548	13.59%	1.191%
47	14.385	1873	1880	1886	rBV	346598	549965	49.31%	4.323%
48	14.578	1910	1913	1916	rVB	3286	3856	0.35%	0.030%
49	14.690	1925	1932	1939	rBV	518917	785730	70.45%	6.176%
50	14.755	1939	1943	1947	rVB2	3987	6317	0.57%	0.050%
51	14.872	1957	1963	1970	rBV2	105074	192185	17.23%	1.511%
52	14.937	1970	1974	1981	rVV4	19820	44741	4.01%	0.352%
53	15.025	1985	1989	1996	rVV6	6125	11635	1.04%	0.091%
54	15.090	1996	2000	2007	rVB7	8186	13773	1.23%	0.108%
55	15.477	2060	2066	2071	rBV5	4065	8918	0.80%	0.070%
56	15.524	2071	2074	2078	rVV3	2325	3346	0.30%	0.026%
57	15.683	2094	2101	2107	rBV	498334	747382	67.01%	5.875%
58	15.795	2113	2120	2127	rBV2	70511	116004	10.40%	0.912%
59	15.859	2129	2131	2134	rVV2	2885	2724	0.24%	0.021%
60	15.924	2138	2142	2148	rVB3	7147	11830	1.06%	0.093%
61	16.018	2156	2158	2160	rBV3	2981	2183	0.20%	0.017%
62	16.253	2197	2198	2205	rVB3	3388	4199	0.38%	0.033%
63	16.365	2215	2217	2218	rBV	2489	2084	0.19%	0.016%
64	16.388	2218	2221	2222	rBV2	4772	4382	0.39%	0.034%
65	16.453	2229	2232	2233	rBV	2589	2722	0.24%	0.021%
66	16.470	2233	2235	2238	rVB3	4691	3707	0.33%	0.029%
67	16.500	2238	2240	2243	rVB2	2584	2413	0.22%	0.019%
68	16.623	2257	2261	2265	rBV4	1995	4241	0.38%	0.033%
69	16.699	2273	2274	2276	rBV2	3117	2379	0.21%	0.019%
70	16.729	2278	2279	2281	rBV2	3638	2150	0.19%	0.017%
71	16.817	2292	2294	2297	rBV3	4440	4059	0.36%	0.032%

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG102920\
 Data File : BG047118.D
 Acq On : 29 Oct 2020 13:43
 Operator : CG/JU
 Sample : L4499-06
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 EW2B7

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_G\METHODS\SOM-EPA-BG102220MA.M
 Title : SVOA CALIBRATION

72	16.911	2307	2310	2311	rVB3	3205	2569	0.23%	0.020%
73	16.964	2316	2319	2323	rVB6	6309	8660	0.78%	0.068%
74	17.005	2323	2326	2327	rBV2	2672	3554	0.32%	0.028%
75	17.023	2327	2329	2330	rBV2	3147	2971	0.27%	0.023%
76	17.076	2335	2338	2339	rBV3	4523	4233	0.38%	0.033%
77	17.181	2353	2356	2360	rVB2	6938	7993	0.72%	0.063%
78	17.316	2375	2379	2386	rBV6	16040	26655	2.39%	0.210%
79	17.381	2388	2390	2393	rVV4	8103	11048	0.99%	0.087%
80	17.434	2393	2399	2404	rVV	666025	1017351	91.22%	7.997%
81	17.475	2404	2406	2411	rVV	74949	94387	8.46%	0.742%
82	17.534	2411	2416	2425	rVB	546447	806945	72.35%	6.343%
83	17.898	2474	2478	2483	rVB2	24041	39404	3.53%	0.310%
84	18.074	2503	2508	2514	rBV2	124998	182385	16.35%	1.434%
85	18.127	2514	2517	2520	rBV5	5134	5821	0.52%	0.046%
86	18.198	2526	2529	2534	rBV7	12605	16871	1.51%	0.133%
87	18.256	2535	2539	2544	rBV2	53269	74112	6.65%	0.583%
88	18.315	2544	2549	2554	rBV2	188065	278637	24.98%	2.190%
89	18.803	2627	2632	2635	rBV6	17826	34971	3.14%	0.275%
90	18.873	2639	2644	2650	rVB	168827	229010	20.53%	1.800%
91	18.985	2659	2663	2669	rVB2	92036	118884	10.66%	0.934%
92	19.114	2681	2685	2689	rBV3	47404	63548	5.70%	0.500%
93	19.208	2696	2701	2705	rBV	131740	180483	16.18%	1.419%
94	19.414	2731	2736	2741	rBV	362173	445517	39.95%	3.502%
95	19.484	2743	2748	2755	rVB	182692	264111	23.68%	2.076%
96	19.578	2761	2764	2770	rBV5	62448	91391	8.19%	0.718%
97	19.690	2779	2783	2788	rBV	203489	266213	23.87%	2.093%
98	19.819	2799	2805	2815	rBV2	631355	1115280	100.00%	8.766%
99	21.705	3121	3126	3131	rVB	615523	945278	84.76%	7.430%
100	21.864	3149	3153	3156	rBV	239543	354479	31.78%	2.786%

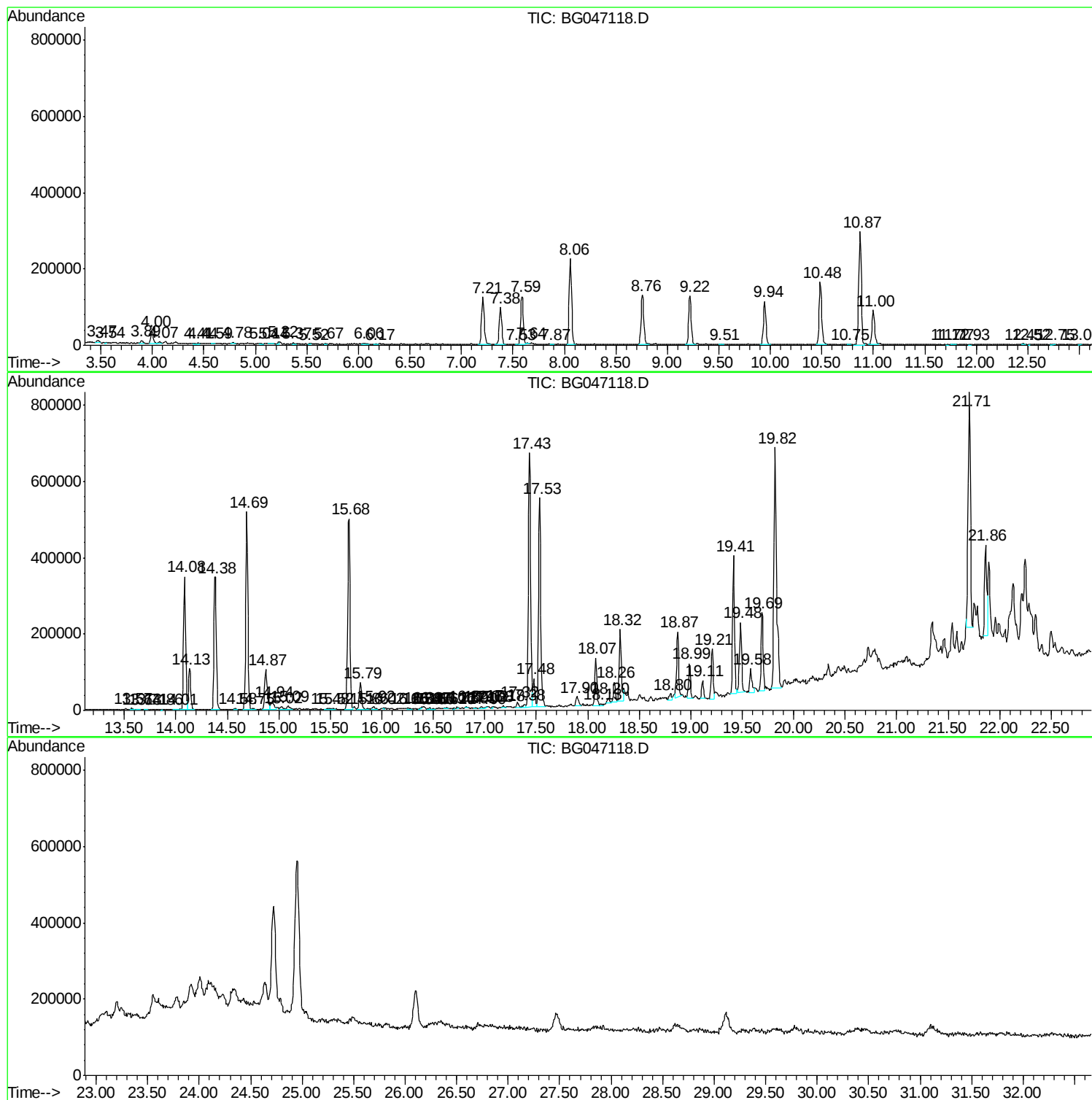
Sum of corrected areas: 12722101

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG102920\
Data File : BG047118.D
Acq On : 29 Oct 2020 13:43
Operator : CG/JU
Sample : L4499-06
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
EW2B7

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG102220MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG102920\
Data File : BG047118.D
Acq On : 29 Oct 2020 13:43
Operator : CG/JU
Sample : L4499-06
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
EW2B7

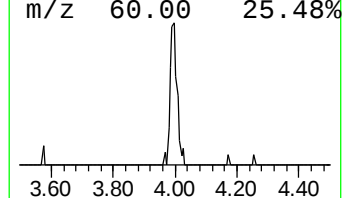
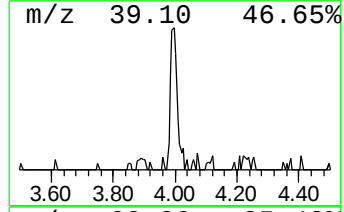
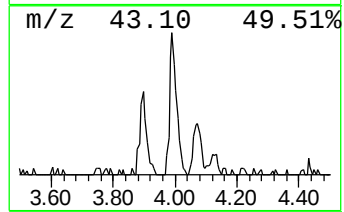
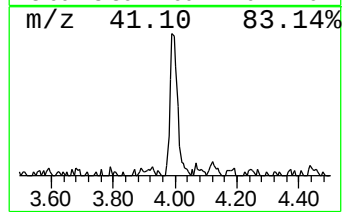
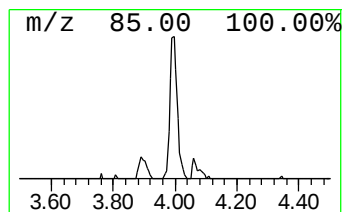
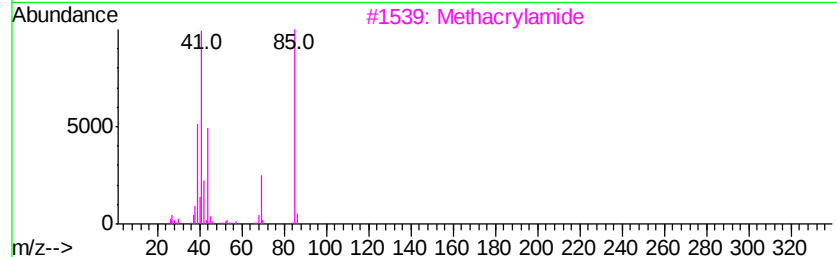
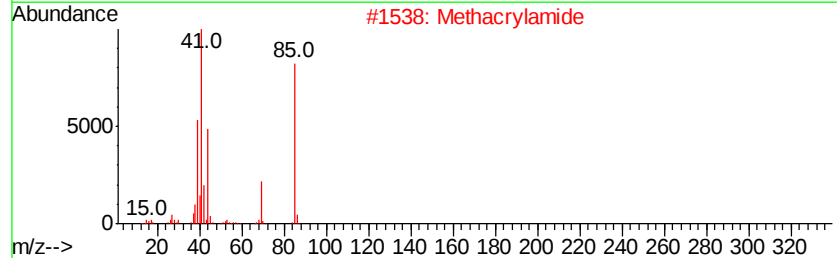
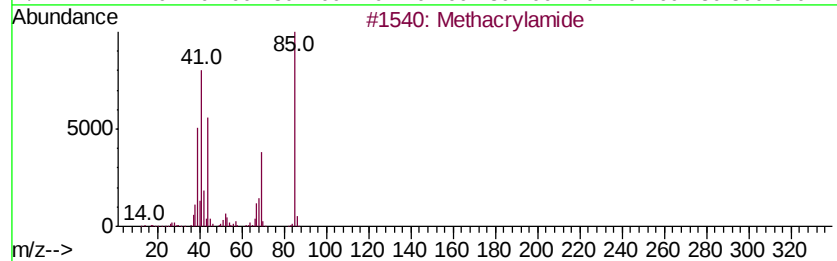
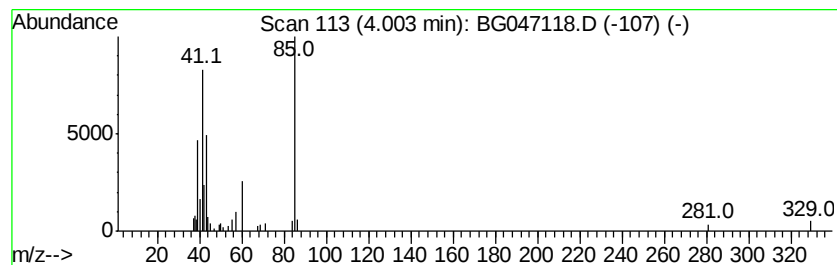
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG102220MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 Methacrylamide Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.00	2.87 ng/ul	54773	1,4-Dichlorobenzene-d4	8.06

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Methacrylamide	85	C4H7NO	000079-39-0	59
2			Methacrylamide	85	C4H7NO	000079-39-0	53
3			Methacrylamide	85	C4H7NO	000079-39-0	53
4			2-(Chloromethyl)tetrahydropyran	134	C6H11ClO	018420-41-2	43
5			2-(Bromomethyl)acrylic acid	164	C4H5BrO2	072707-66-5	36



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG102920\
Data File : BG047118.D
Acq On : 29 Oct 2020 13:43
Operator : CG/JU
Sample : L4499-06
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
EW2B7

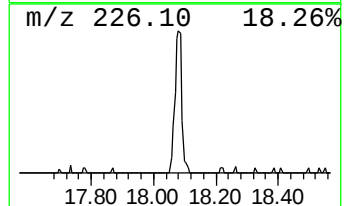
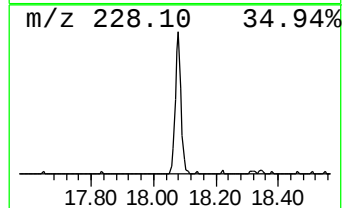
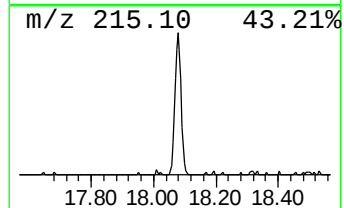
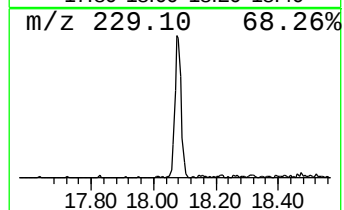
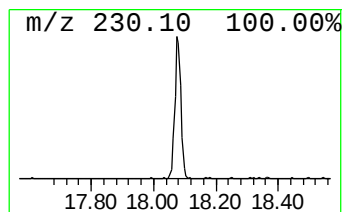
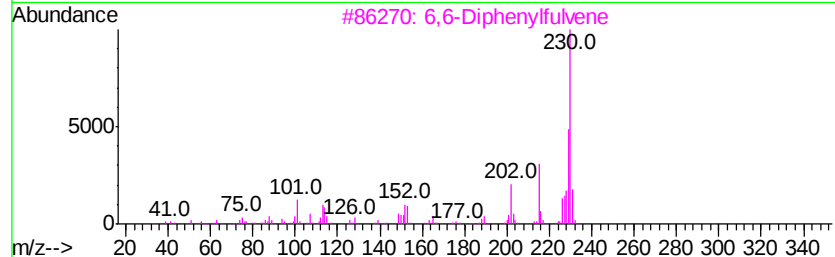
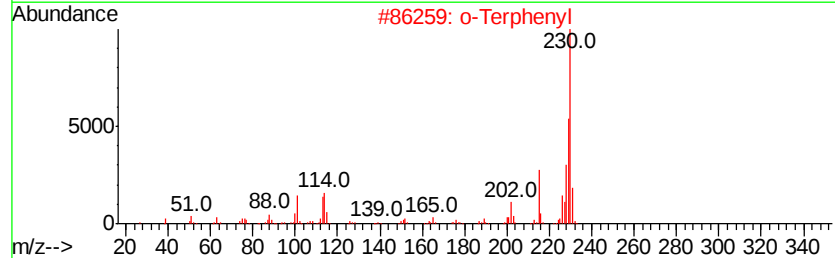
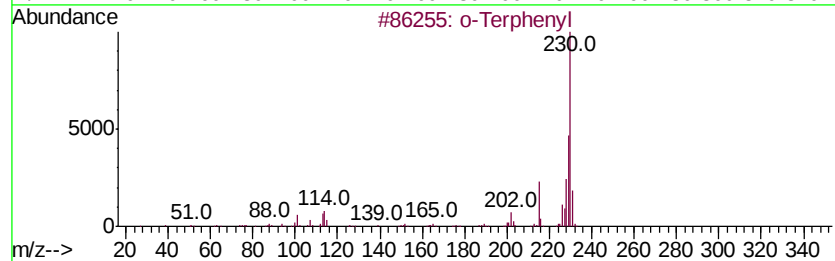
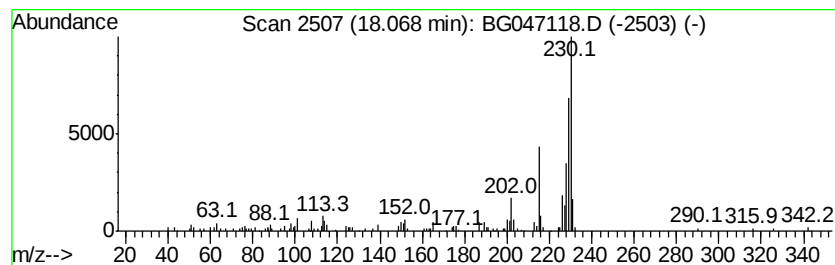
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG102220MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 o-Terphenyl Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.07	3.59 ng/ul	182385	Phenanthrene-d10	17.43

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	o-Terphenyl	230	C18H14	000084-15-1	93
2		o-Terphenyl	230	C18H14	000084-15-1	91
3		6,6-Diphenylfulvene	230	C18H14	002175-90-8	90
4		o-Terphenyl	230	C18H14	000084-15-1	87
5		6,6-Diphenylfulvene	230	C18H14	002175-90-8	68



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG102920\
Data File : BG047118.D
Acq On : 29 Oct 2020 13:43
Operator : CG/JU
Sample : L4499-06
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
EW2B7

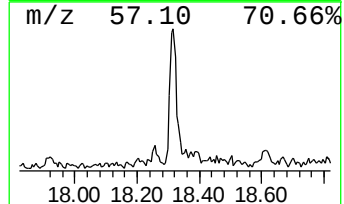
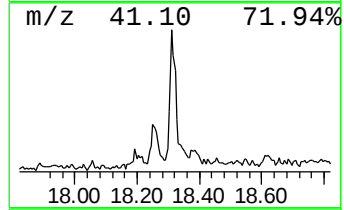
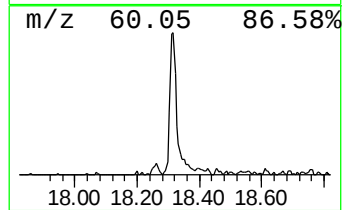
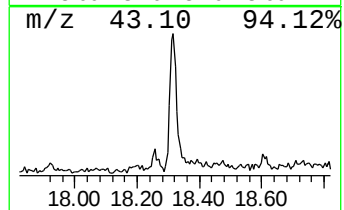
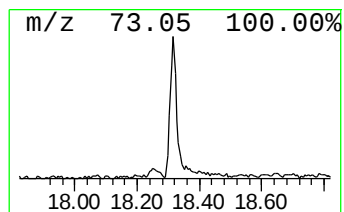
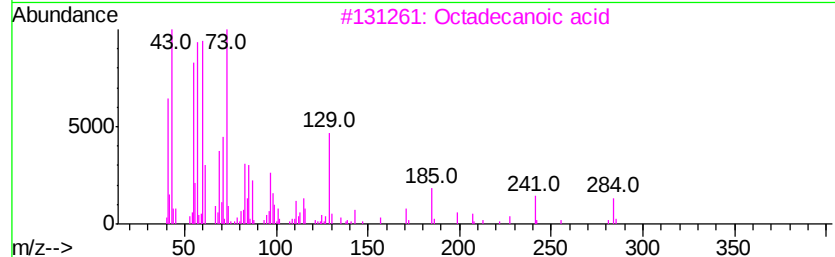
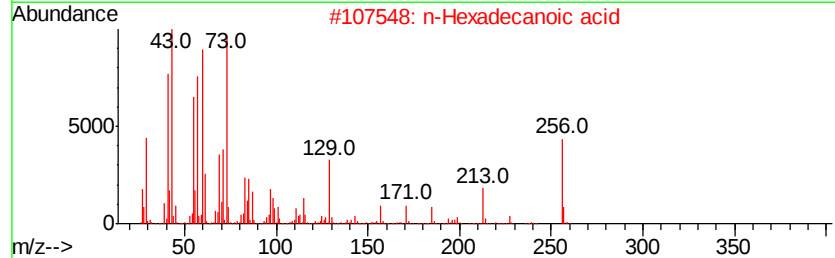
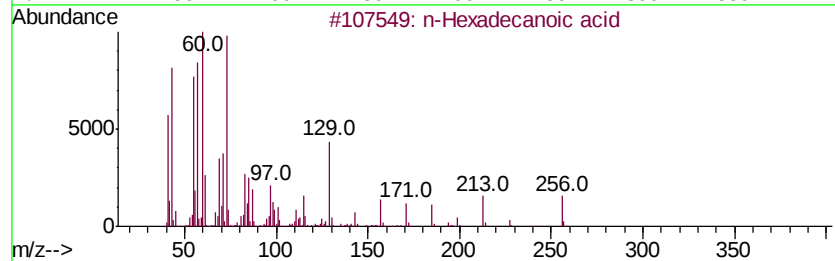
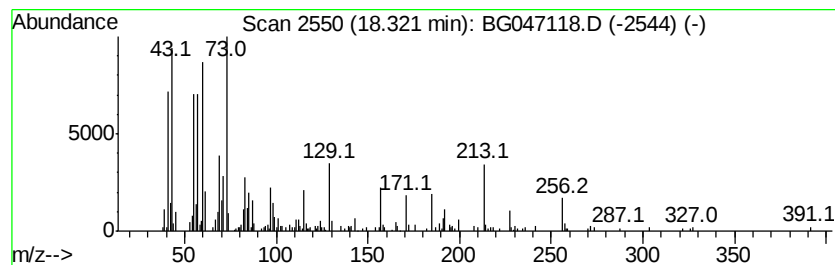
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG102220MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 n-Hexadecanoic acid Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.32	5.48 ng/ul	278637	Phenanthrene-d10	17.43

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98	
2	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	97	
3	Octadecanoic acid	284	C18H36O2	000057-11-4	83	
4	Tetradecanoic acid	228	C14H28O2	000544-63-8	80	
5	Tridecanoic acid	214	C13H26O2	000638-53-9	76	



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG102920\
Data File : BG047118.D
Acq On : 29 Oct 2020 13:43
Operator : CG/JU
Sample : L4499-06
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
EW2B7

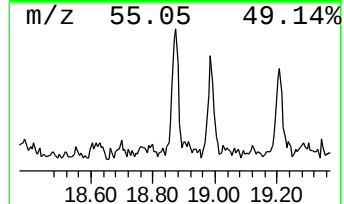
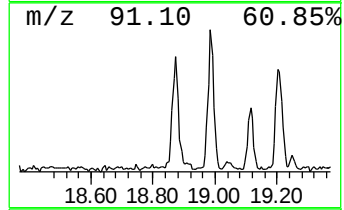
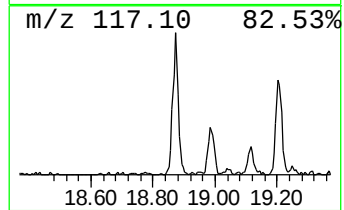
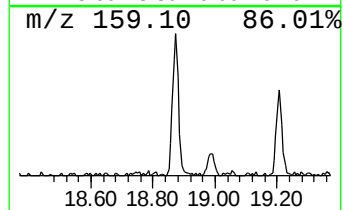
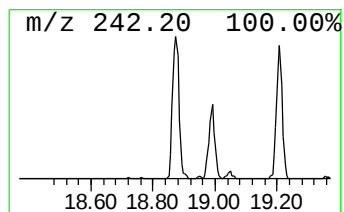
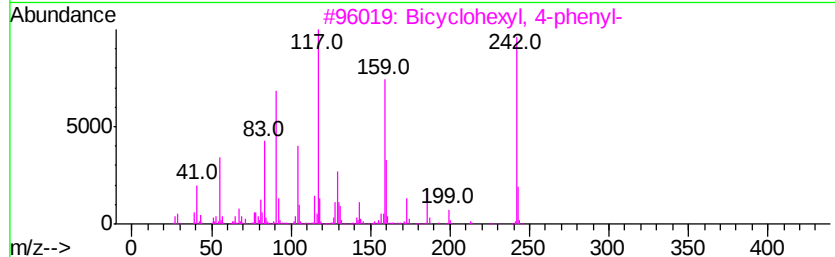
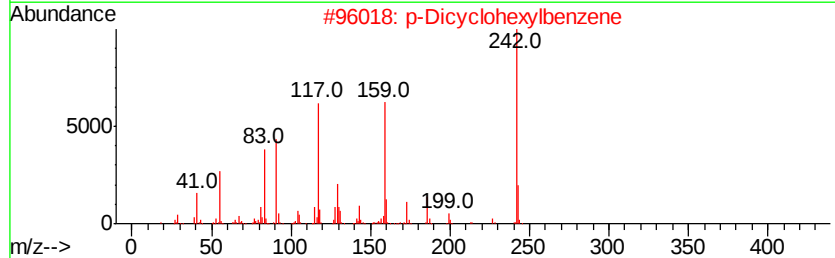
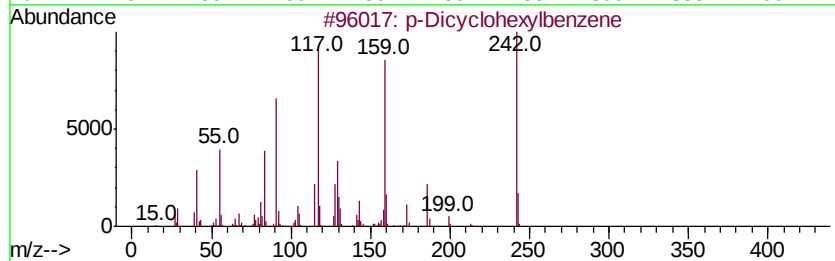
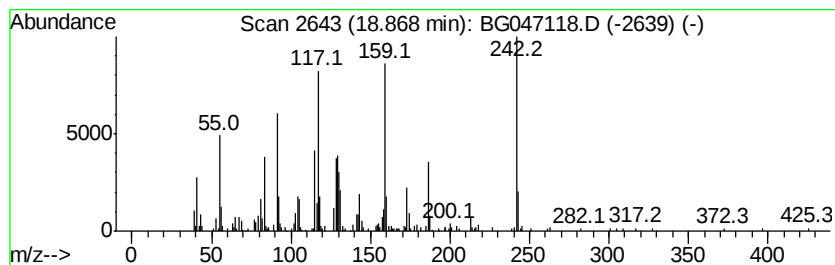
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG102220MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 p-Dicyclohexylbenzene Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.87	4.50 ng/ul	229010	Phenanthrene-d10	17.43

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		p-Dicvclohexylbenzene	242	C18H26	001087-02-1	96
2		p-Dicvclohexylbenzene	242	C18H26	001087-02-1	83
3		Bicvclohexyl, 4-phenvl-	242	C18H26	020273-27-2	76
4		1,1-Diethyl-1,2,3,4-tetrahydrona...	188	C14H20	002938-66-1	35
5		2-Naphthalenol, 5-amino-	159	C10H9NO	000086-97-5	25



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG102920\
Data File : BG047118.D
Acq On : 29 Oct 2020 13:43
Operator : CG/JU
Sample : L4499-06
Misc :
ALS Vial : 9 Sample Multiplier: 1

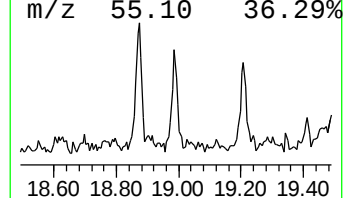
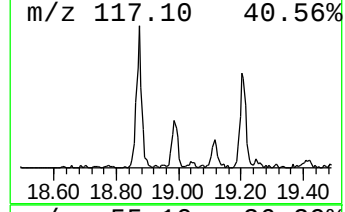
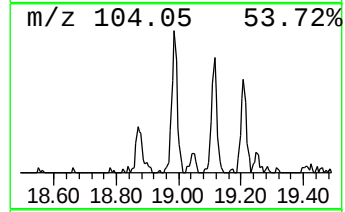
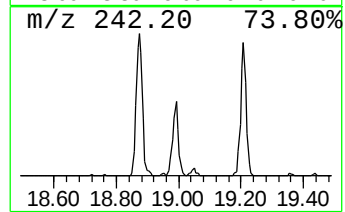
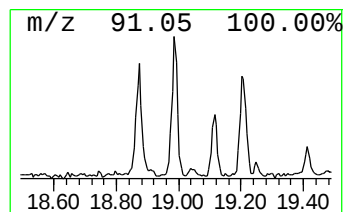
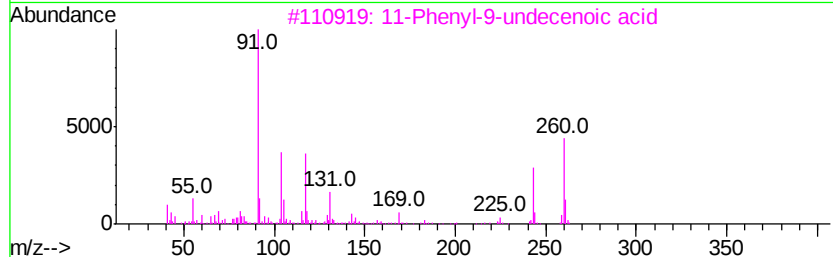
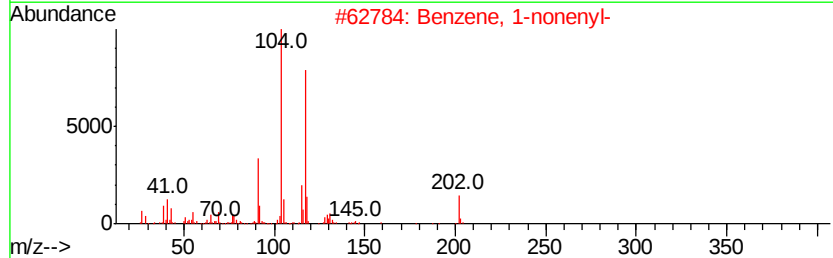
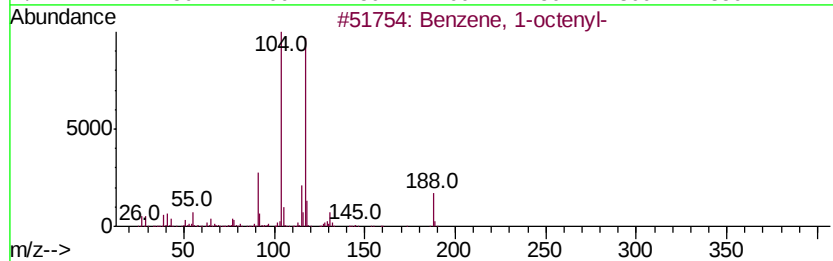
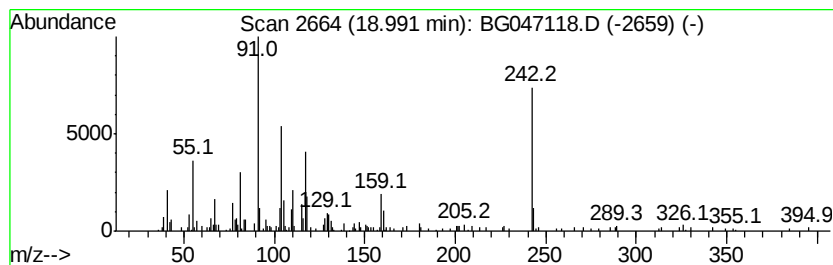
Instrument :
BNA_G
ClientSampleId :
EW2B7

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG102220MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 unknown-01 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.		
18.99	2.34 ng/ul	118884	Phenanthrene-d10	17.43		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Benzene, 1-octenyl-	188	C14H20	029518-72-7	38
2		Benzene, 1-nonenyl-	202	C15H22	034426-61-4	38
3		11-Phenyl-9-undecenoic acid	260	C17H24O2	1000131-48-9	16
4		Benzaldehyde, 2-hydroxy-6-methyl...	242	C15H14O3	034883-10-8	10
5		6-Dimethylaminomethyltricyclo[8....	281	C19H23NO	1000306-67-8	10



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG102920\
Data File : BG047118.D
Acq On : 29 Oct 2020 13:43
Operator : CG/JU
Sample : L4499-06
Misc :
ALS Vial : 9 Sample Multiplier: 1

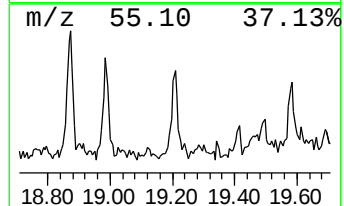
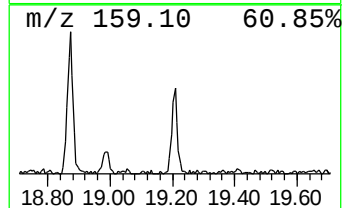
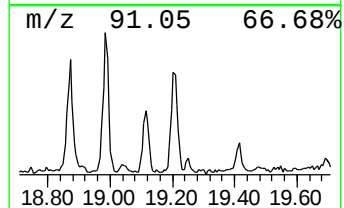
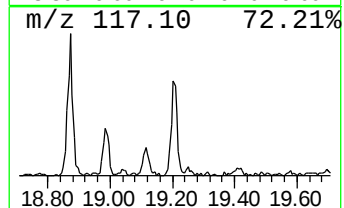
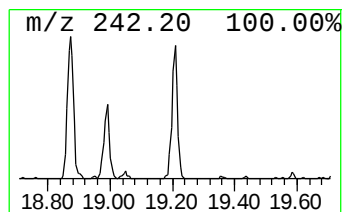
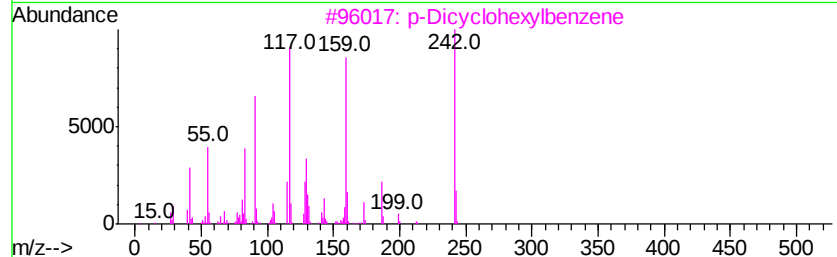
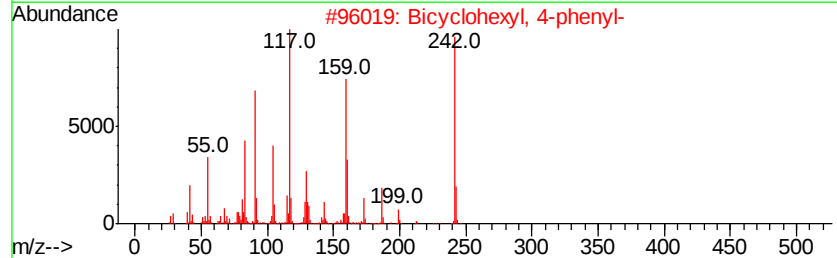
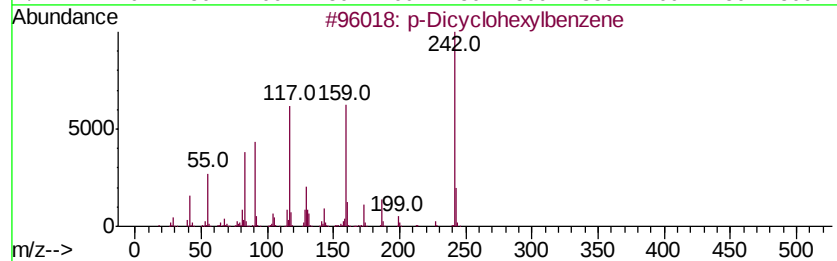
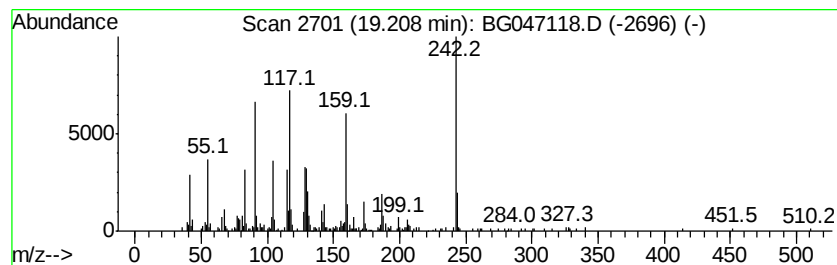
Instrument :
BNA_G
ClientSampleId :
EW2B7

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG102220MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 Bicyclohexyl, 4-phenyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.		
19.21	3.55 ng/ul	180483	Phenanthrene-d10	17.43		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		p-Dicvclohexylbenzene	242	C18H26	001087-02-1	96
2		Bicvclohexyl, 4-phenyl-	242	C18H26	020273-27-2	93
3		p-Dicvclohexylbenzene	242	C18H26	001087-02-1	91
4		9H-Tribenzo[a,c,e)cycloheptene	242	C19H14	000213-10-5	30
5		1,3-Diamino-5,6-dihydro-8-methox...	242	C13H14N4O	037436-50-3	27



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG102920\
Data File : BG047118.D
Acq On : 29 Oct 2020 13:43
Operator : CG/JU
Sample : L4499-06
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
EW2B7

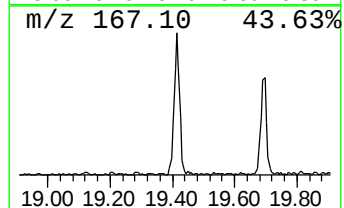
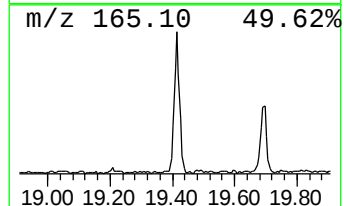
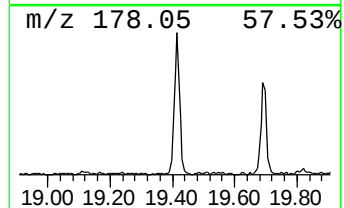
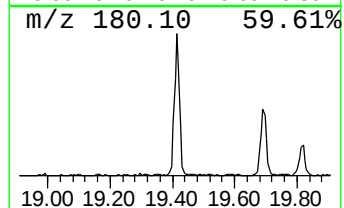
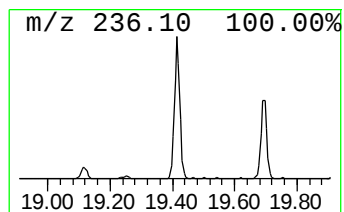
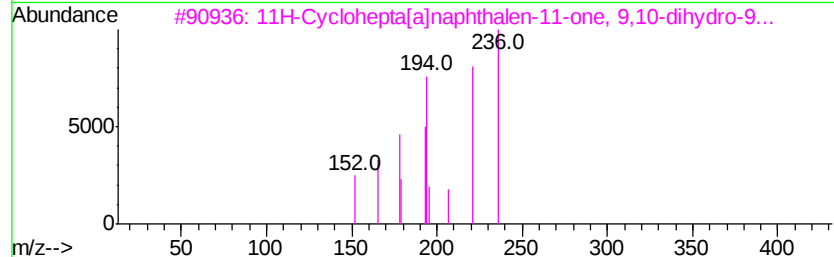
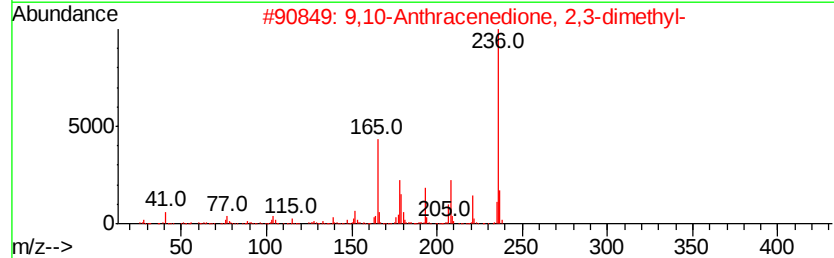
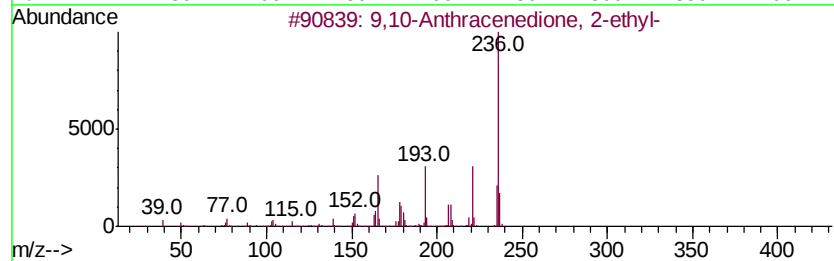
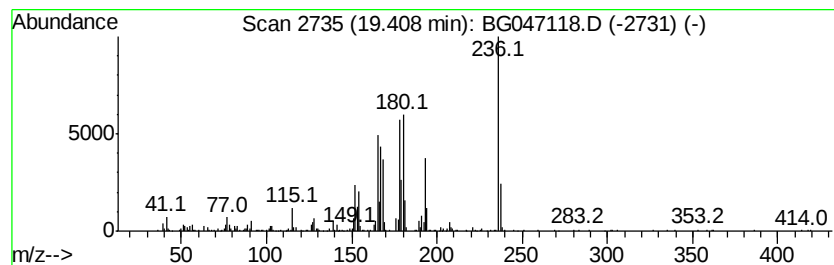
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG102220MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 7 unknown-02 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.41	8.76 ng/ul	445517	Phenanthrene-d10	17.43

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	9,10-Anthracenedione, 2-ethyl-	236	C16H12O2	000084-51-5	46
2		9,10-Anthracenedione, 2,3-dimethyl-	236	C16H12O2	006531-35-7	38
3		11H-Cyclohepta[fa]naphthalen-11-one...	236	C17H16O	058111-76-5	38
4		9,10-Anthracenedione, 2-ethyl-	236	C16H12O2	000084-51-5	38
5		2,4-Diethoxy-5-methoxy-1-(2-prop...	236	C14H20O3	1000122-72-0	35



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG102920\
Data File : BG047118.D
Acq On : 29 Oct 2020 13:43
Operator : CG/JU
Sample : L4499-06
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
EW2B7

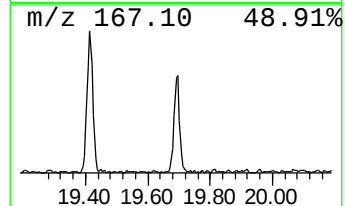
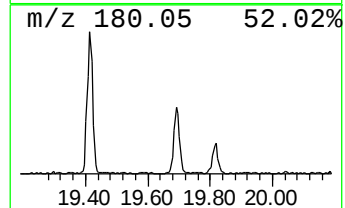
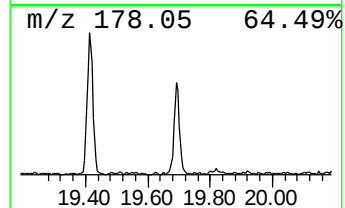
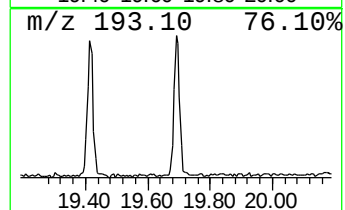
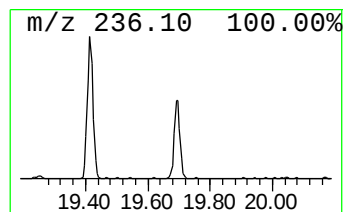
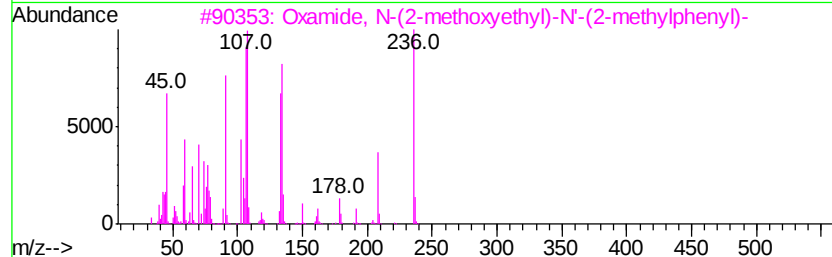
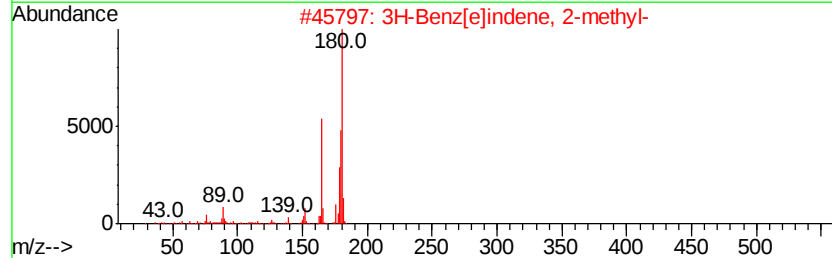
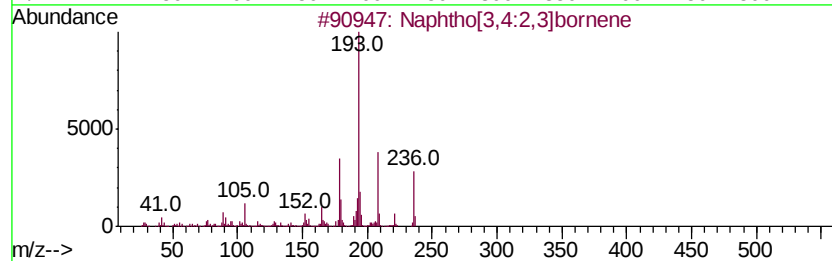
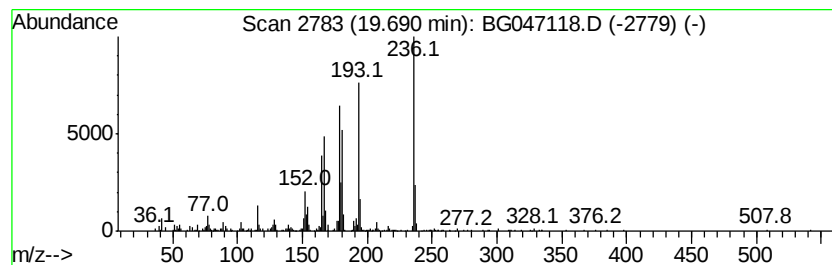
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG102220MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 8 Naphtho[3,4:2,3]bornene Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.69	5.63 ng/ul	266213	Chrysene-d12	21.71

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphtho[3,4:2,3]bornene	236	C18H20	1000210-68-3	76
2		3H-Benz[e]indene, 2-methyl-	180	C14H12	150096-60-9	56
3		Oxamide, N-(2-methoxyethyl)-N'-(...	236	C12H16N2O3	1000277-54-5	53
4		Carbamazepine	236	C15H12N2O	000298-46-4	43
5		Ethylene, 1,1-diphenyl-	180	C14H12	000530-48-3	25



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG102920\
Data File : BG047118.D
Acq On : 29 Oct 2020 13:43
Operator : CG/JU
Sample : L4499-06
Misc :
ALS Vial : 9 Sample Multiplier: 1

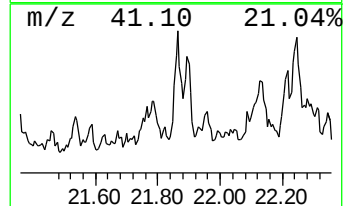
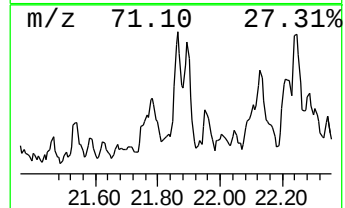
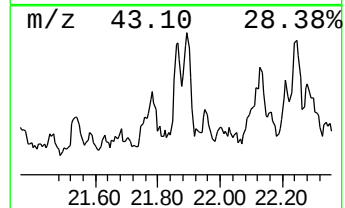
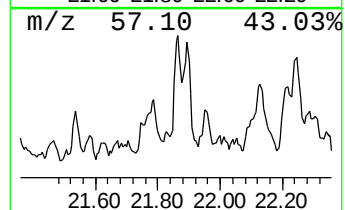
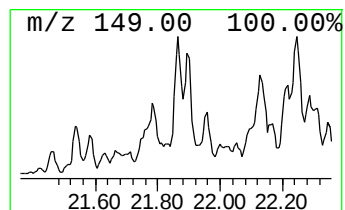
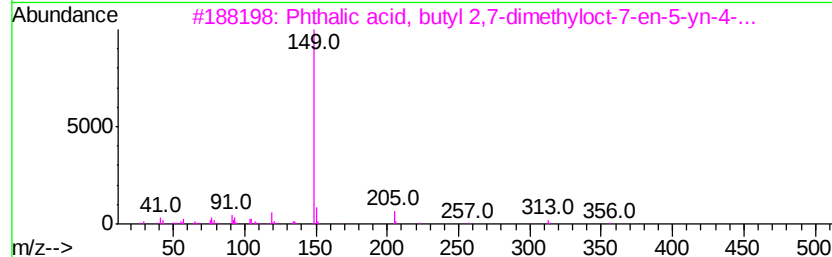
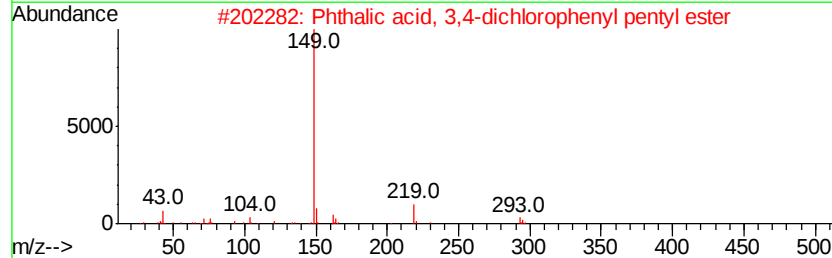
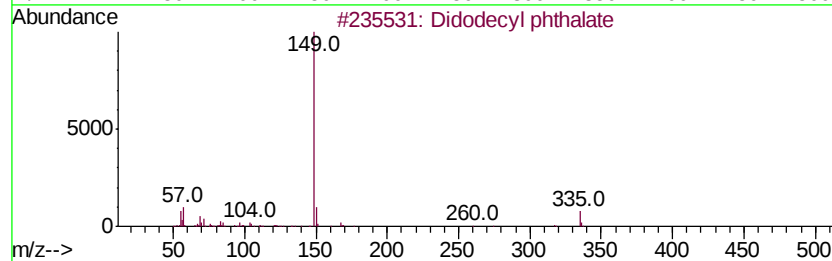
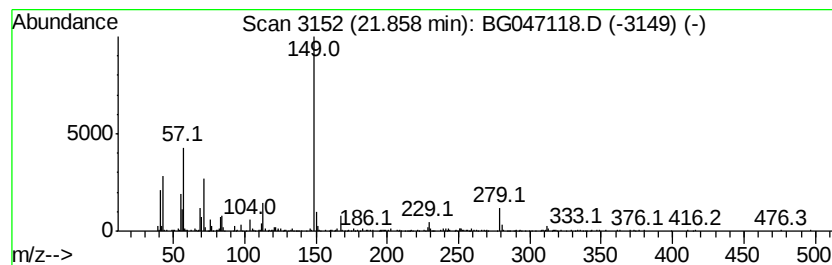
Instrument :
BNA_G
ClientSampleId :
EW2B7

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG102220MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 9 unknown-03 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.		
21.86	7.50 ng/ul	354479	Chrysene-d12	21.71		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Didodecyl phthalate	502	C32H54O4	002432-90-8	47
2		Phthalic acid, 3,4-dichloropheny...	380	C19H18Cl2O4	1000315-27-4	47
3		Phthalic acid, butyl 2,7-dimethv...	356	C22H28O4	1000315-49-0	46
4		Phthalic acid, ethyl pentyl ester	264	C15H20O4	1000308-93-6	46
5		Diisooctyl phthalate	390	C24H38O4	000131-20-4	43



Data Path : Z:\SVOASRV\HPCHEM1\BNA_G\DATA\BG102920\
Data File : BG047118.D
Acq On : 29 Oct 2020 13:43
Operator : CG/JU
Sample : L4499-06
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
EW2B7

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG102220MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Methacrylamide	4.00	2.9	ng/ul	54773	1	8.06	381563	20.0
o-Terphenyl	18.07	3.6	ng/ul	182385	4	17.43	1017350	20.0
n-Hexadecanoic acid	18.32	5.5	ng/ul	278637	4	17.43	1017350	20.0
p-Dicyclohexylben...	18.87	4.5	ng/ul	229010	4	17.43	1017350	20.0
unknown-01	18.99	2.3	ng/ul	118884	4	17.43	1017350	20.0
Bicyclohexyl, 4-p...	19.21	3.5	ng/ul	180483	4	17.43	1017350	20.0
unknown-02	19.41	8.8	ng/ul	445517	4	17.43	1017350	20.0
Naphtho[3,4:2,3]b...	19.69	5.6	ng/ul	266213	5	21.71	945278	20.0
unknown-03	21.86	7.5	ng/ul	354479	5	21.71	945278	20.0