

Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM052620\
 Data File : BM026126.D
 Acq On : 26 May 2020 23:05
 Operator : MA/SJ
 Sample : L2756-09
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 ETKX1

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_M\METHODS\SOM-EPA-BM051320MA.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	3.028	20	24	27	rBV	30069	41585	0.54%	0.065%
2	3.063	27	30	38	rVB	52745	71888	0.93%	0.112%
3	3.763	144	149	161	rVB	554068	752164	9.74%	1.170%
4	4.275	230	236	247	rVB	305515	431952	5.59%	0.672%
5	5.045	362	367	376	rVB	134920	188943	2.45%	0.294%
6	5.193	383	392	401	rBV	274271	496865	6.43%	0.773%
7	5.510	440	446	448	rBV	513483	747994	9.69%	1.164%
8	6.004	525	530	536	rBV	33121	48289	0.63%	0.075%
9	6.598	625	631	635	rBV	30785	43621	0.56%	0.068%
10	6.681	635	645	650	rVV2	167150	310722	4.02%	0.483%
11	6.728	650	653	660	rVB	73925	102348	1.33%	0.159%
12	6.834	665	671	678	rBV	539142	851519	11.03%	1.325%
13	6.887	678	680	684	rVV	48917	68405	0.89%	0.106%
14	6.928	684	687	692	rVB3	28282	42775	0.55%	0.067%
15	7.028	697	704	711	rVV	593314	911713	11.81%	1.419%
16	7.145	718	724	733	rVV3	245447	437149	5.66%	0.680%
17	7.234	733	739	746	rVB	45047	69458	0.90%	0.108%
18	7.487	775	782	788	rBV	651612	989909	12.82%	1.540%
19	7.604	793	802	807	rVV	105172	170953	2.21%	0.266%
20	7.669	809	813	819	rVB	45127	67583	0.88%	0.105%
21	7.857	838	845	850	rBV	41063	63322	0.82%	0.099%
22	8.016	862	872	884	rBV	1367701	2110862	27.33%	3.284%
23	8.204	896	904	912	rVV	436096	667906	8.65%	1.039%
24	8.628	970	976	990	rVV	695908	1190013	15.41%	1.852%
25	8.751	992	997	1005	rBV3	32433	61662	0.80%	0.096%
26	9.104	1052	1057	1062	rVB2	48963	73101	0.95%	0.114%
27	9.163	1062	1067	1073	rVB	58572	91788	1.19%	0.143%
28	9.239	1073	1080	1085	rBV2	22276	34918	0.45%	0.054%
29	9.345	1090	1098	1109	rBV	571555	911010	11.80%	1.417%
30	9.686	1146	1156	1163	rBV3	212110	508180	6.58%	0.791%
31	9.757	1163	1168	1172	rVV	148878	254057	3.29%	0.395%
32	9.792	1172	1174	1182	rVV	72961	112702	1.46%	0.175%
33	9.880	1182	1189	1198	rVV	811373	1299627	16.83%	2.022%
34	9.957	1198	1202	1212	rVB4	20982	46482	0.60%	0.072%

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Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM051320MA.M
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35	10.245	1242	1251	1255	rVV	833664	1441492	18.67%	2.243%
36	10.298	1255	1260	1269	rVV	4864439	7722208	100.00%	12.015%
37	10.410	1269	1279	1287	rVV2	200829	423609	5.49%	0.659%
38	11.445	1450	1455	1461	rBV	24138	39998	0.52%	0.062%
39	11.586	1468	1479	1484	rVB4	22399	47551	0.62%	0.074%
40	11.810	1513	1517	1521	rBV	34461	50176	0.65%	0.078%
41	11.922	1530	1536	1542	rVV	112828	179233	2.32%	0.279%
42	12.145	1562	1574	1583	rBV	415013	694749	9.00%	1.081%
43	12.957	1706	1712	1723	rBV	428029	629090	8.15%	0.979%
44	13.098	1730	1736	1740	rBV2	24201	44480	0.58%	0.069%
45	13.180	1746	1750	1755	rVB2	36072	58136	0.75%	0.090%
46	13.298	1758	1770	1778	rBV7	31565	78027	1.01%	0.121%
47	13.451	1786	1796	1801	rBV	41589	75383	0.98%	0.117%
48	13.557	1807	1814	1819	rVV	1552850	2199472	28.48%	3.422%
49	13.598	1819	1821	1831	rVB	132281	178418	2.31%	0.278%
50	13.686	1831	1836	1840	rBV2	41627	61965	0.80%	0.096%
51	13.821	1852	1859	1861	rBV	1627519	2507831	32.48%	3.902%
52	13.845	1861	1863	1873	rVB	1166836	1504086	19.48%	2.340%
53	14.127	1905	1911	1918	rBV	1209194	1821691	23.59%	2.834%
54	14.192	1918	1922	1926	rVV	123398	175866	2.28%	0.274%
55	14.227	1926	1928	1933	rVB	26911	32793	0.42%	0.051%
56	14.357	1945	1950	1959	rBV	124019	221344	2.87%	0.344%
57	14.439	1961	1964	1972	rVV3	17141	35649	0.46%	0.055%
58	14.533	1976	1980	1988	rVB	26634	45709	0.59%	0.071%
59	14.763	2011	2019	2025	rVV8	12205	33446	0.43%	0.052%
60	15.010	2055	2061	2066	rVB	27484	44909	0.58%	0.070%
61	15.133	2075	2082	2087	rVV	2297197	3164784	40.98%	4.924%
62	15.186	2087	2091	2098	rVV2	75643	129060	1.67%	0.201%
63	15.262	2098	2104	2110	rVV	822118	1091277	14.13%	1.698%
64	15.310	2110	2112	2116	rVV3	30033	41375	0.54%	0.064%
65	15.833	2193	2201	2204	rBV8	20111	41017	0.53%	0.064%
66	16.339	2281	2287	2294	rVB5	22722	40821	0.53%	0.064%
67	16.427	2294	2302	2306	rBV8	13354	32210	0.42%	0.050%
68	16.698	2340	2348	2353	rVV	135307	205846	2.67%	0.320%
69	16.745	2353	2356	2364	rVB3	30038	54151	0.70%	0.084%
70	16.880	2372	2379	2383	rBV	1527483	2102644	27.23%	3.272%
71	16.921	2383	2386	2390	rVV	696904	898477	11.63%	1.398%

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72	16.980	2390	2396	2400	rVV	2636277	3744746	48.49%	5.826%
73	17.009	2400	2401	2410	rVV	317023	324344	4.20%	0.505%
74	17.409	2464	2469	2474	rBV	48939	68930	0.89%	0.107%
75	17.462	2474	2478	2483	rBV2	29974	49208	0.64%	0.077%
76	17.674	2510	2514	2520	rBV4	30594	55943	0.72%	0.087%
77	17.756	2523	2528	2533	rVV	169844	237163	3.07%	0.369%
78	17.804	2533	2536	2540	rVB	56868	70487	0.91%	0.110%
79	17.886	2545	2550	2555	rVV2	36865	58241	0.75%	0.091%
80	17.945	2555	2560	2564	rVV2	227767	373512	4.84%	0.581%
81	17.986	2564	2567	2572	rVB	104047	138931	1.80%	0.216%
82	18.262	2609	2614	2617	rBV	106628	143554	1.86%	0.223%
83	18.298	2617	2620	2626	rVB	118886	158652	2.05%	0.247%
84	18.686	2682	2686	2690	rBV2	41811	63730	0.83%	0.099%
85	18.745	2692	2696	2699	rBV4	25334	40075	0.52%	0.062%
86	18.821	2705	2709	2719	rBV2	56693	99810	1.29%	0.155%
87	18.915	2721	2725	2726	rBV	40318	48006	0.62%	0.075%
88	18.945	2726	2730	2736	rVB	638795	799870	10.36%	1.245%
89	19.098	2752	2756	2761	rVB	52758	70923	0.92%	0.110%
90	19.168	2764	2768	2776	rVB2	33625	73864	0.96%	0.115%
91	19.280	2781	2787	2790	rBV	3477552	4579743	59.31%	7.126%
92	19.309	2790	2792	2801	rVB	886274	958408	12.41%	1.491%
93	19.809	2875	2877	2885	rVB3	44834	58917	0.76%	0.092%
94	19.968	2902	2904	2909	rVB	28548	35378	0.46%	0.055%
95	20.156	2933	2936	2941	rVB2	34876	44156	0.57%	0.069%
96	20.827	3046	3050	3054	rBV	239707	285618	3.70%	0.444%
97	21.080	3087	3093	3097	rBV	1912005	2479069	32.10%	3.857%
98	22.603	3349	3352	3357	rVB2	155458	197504	2.56%	0.307%
99	23.062	3424	3430	3438	rBV	2483013	4214702	54.58%	6.558%
100	23.197	3447	3453	3461	rVB	1358344	2377603	30.79%	3.699%

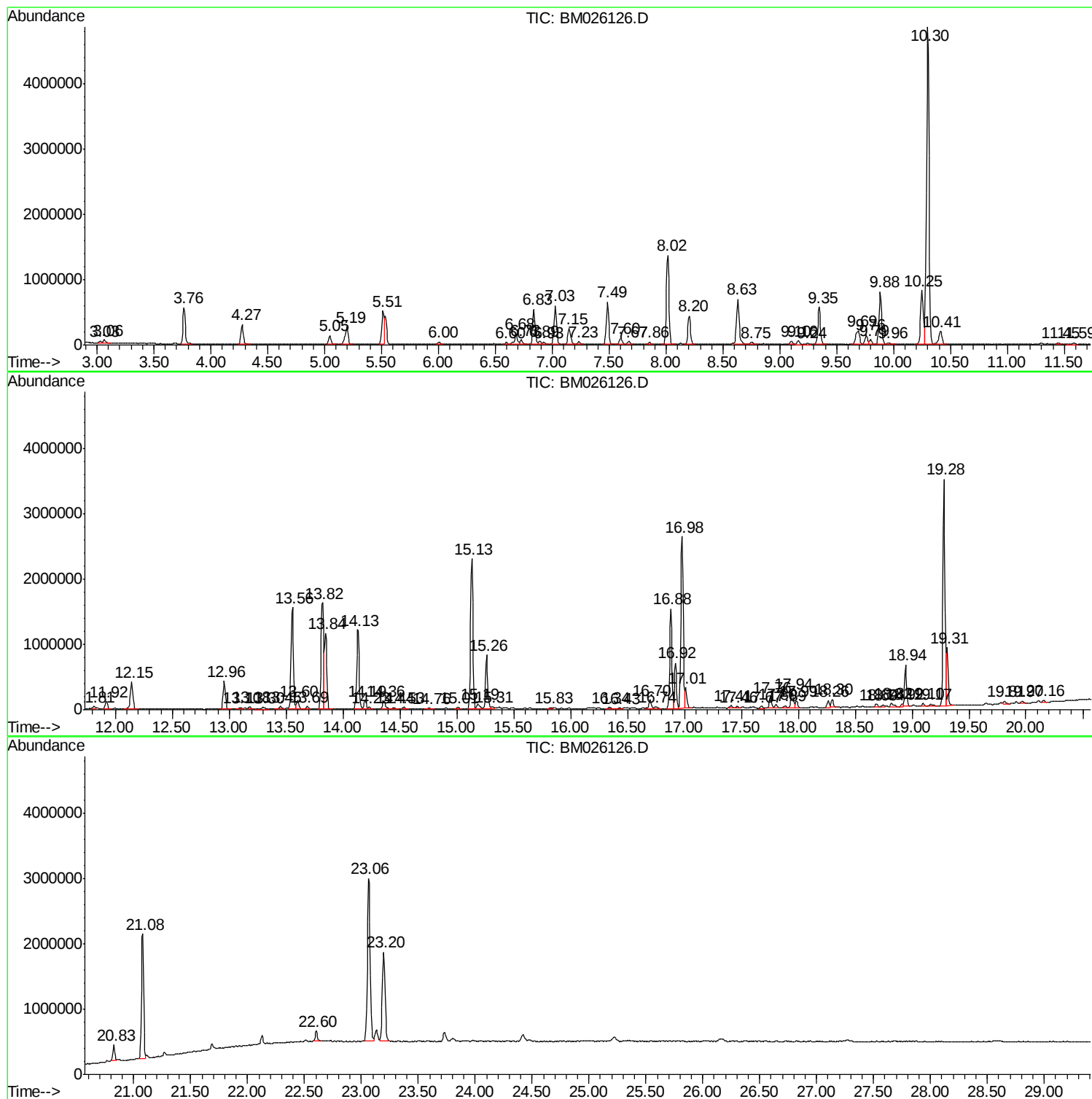
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM051320MA.M
Quant Title : SVOA CALIBRATION

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



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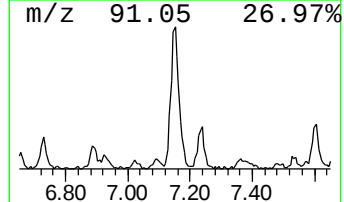
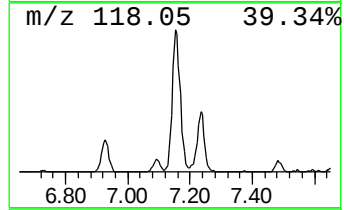
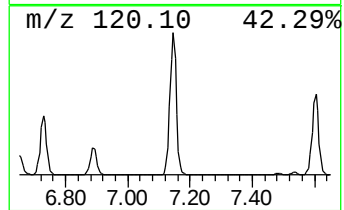
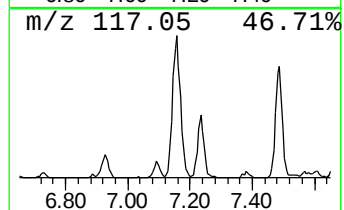
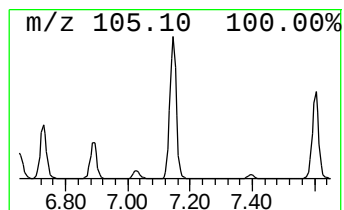
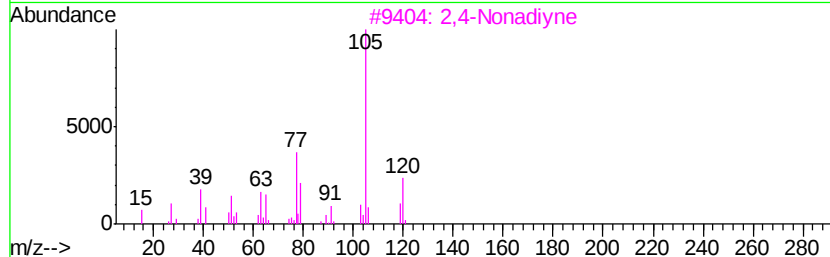
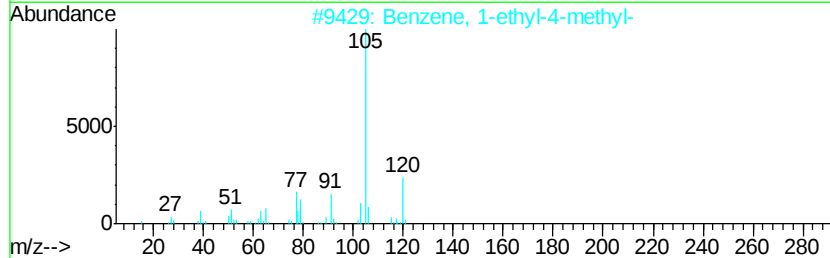
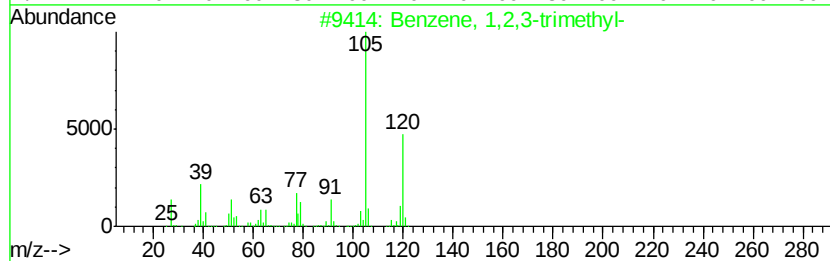
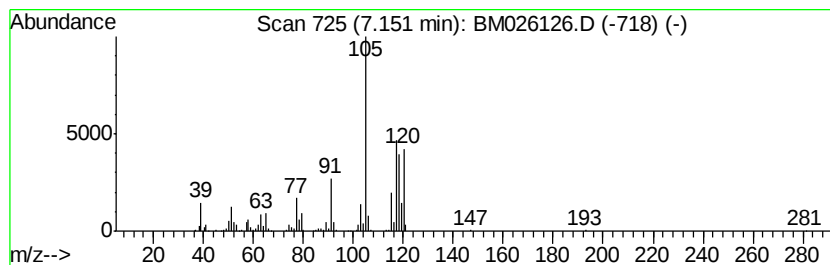
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM051320MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 6 Benzene, 1,2,3-trimethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.15	8.83 ng/ul	437149	1,4-Dichlorobenzene-d4	7.49

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	90
2		Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	87
3		2,4-Nonadiyne	120	C9H12	063621-15-8	64
4		Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	60
5		Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	60



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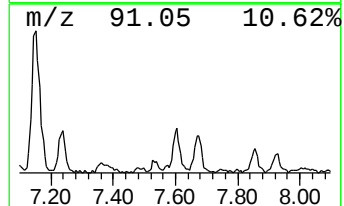
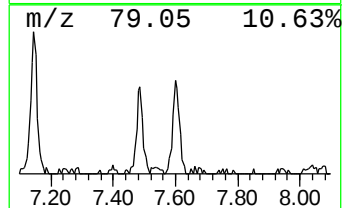
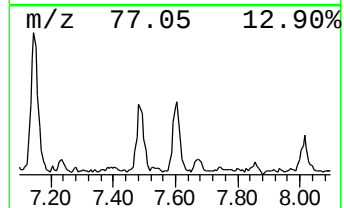
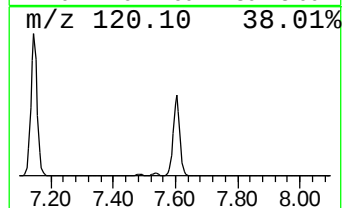
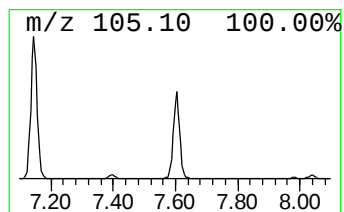
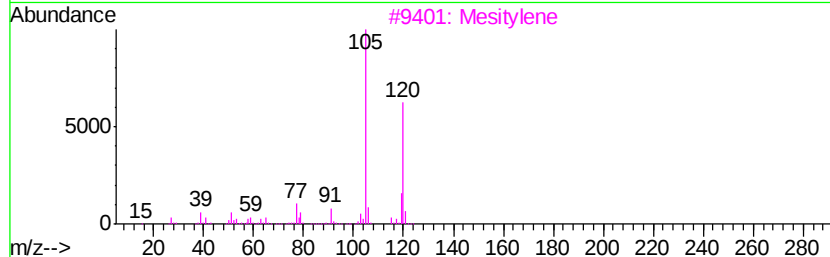
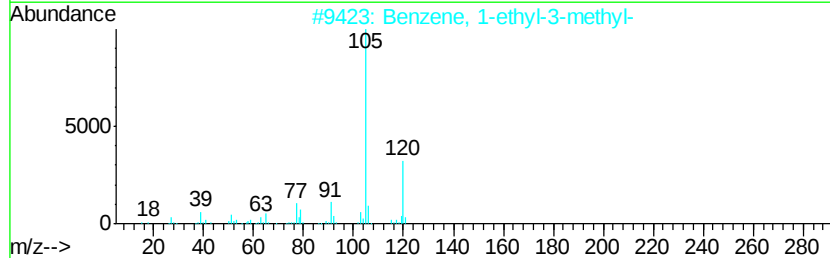
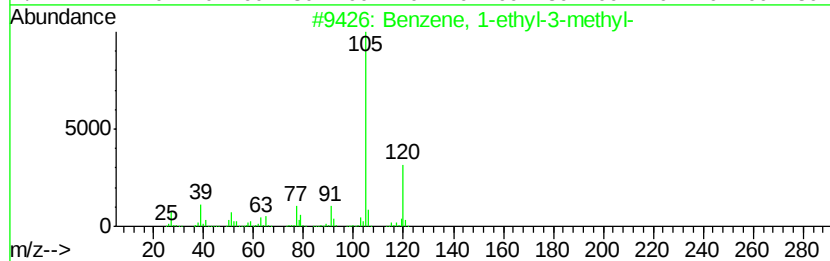
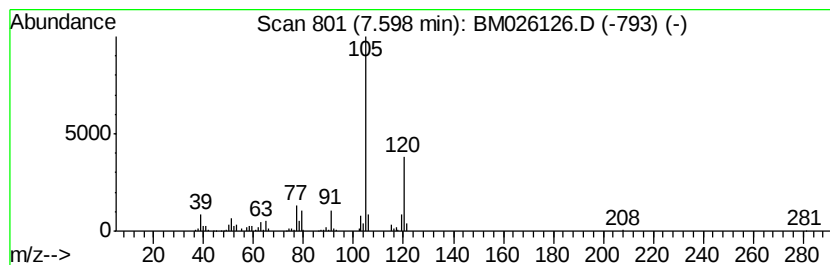
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM051320MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 7 Benzene, 1-ethyl-3-methyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.60	3.45 ng/ul	170953	1,4-Dichlorobenzene-d4	7.49

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91
2		Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91
3		Mesitylene	120	C9H12	000108-67-8	91
4		Mesitylene	120	C9H12	000108-67-8	91
5		Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	91



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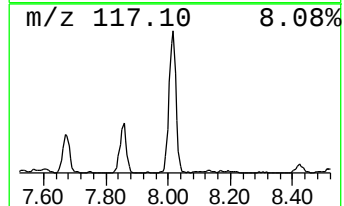
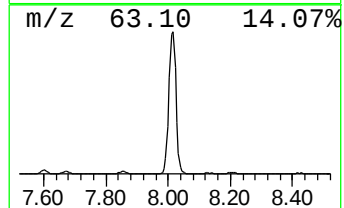
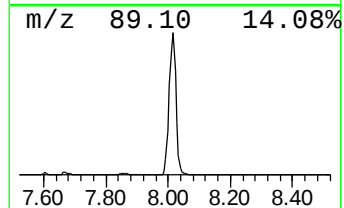
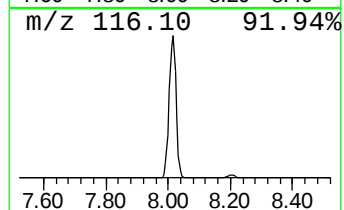
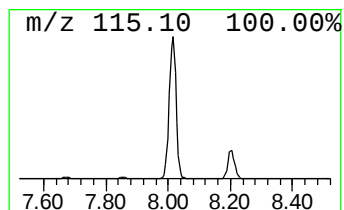
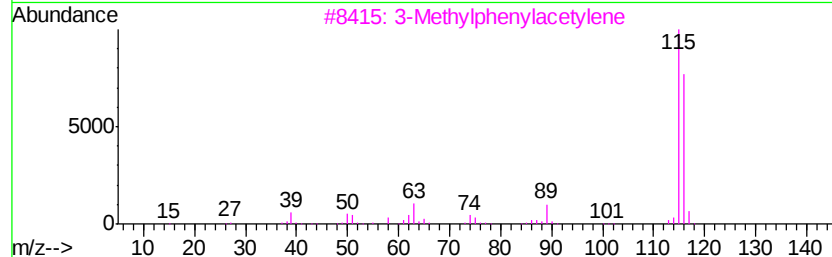
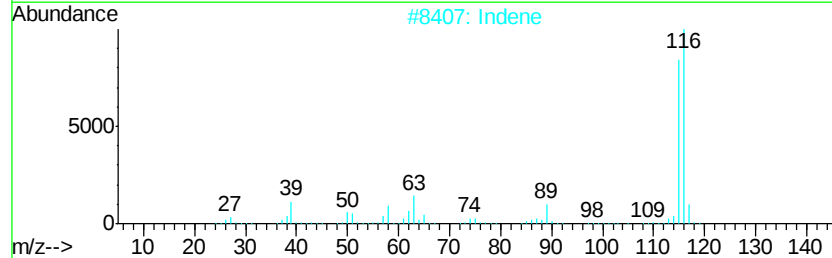
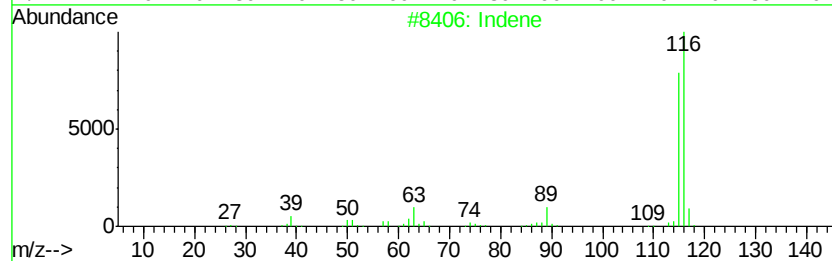
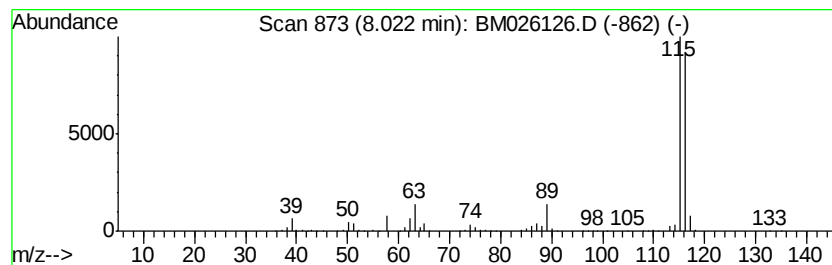
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM051320MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 8 Indene Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.02	42.65 ng/ul	2110860	1,4-Dichlorobenzene-d4	7.49

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Indene	116	C9H8	000095-13-6	94
2		Indene	116	C9H8	000095-13-6	94
3		3-Methylphenylacetylene	116	C9H8	000766-82-5	94
4		Indene	116	C9H8	000095-13-6	93
5		Benzene, 1-propynyl-	116	C9H8	000673-32-5	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM052620\
Data File : BM026126.D
Acq On : 26 May 2020 23:05
Operator : MA/SJ
Sample : L2756-09
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
ETKX1

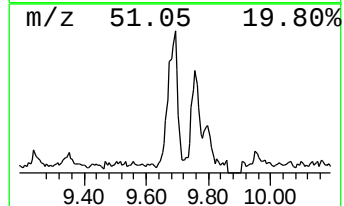
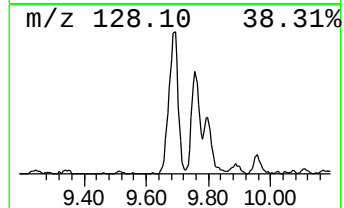
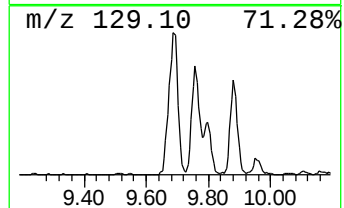
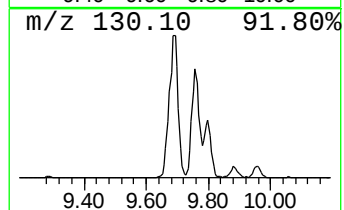
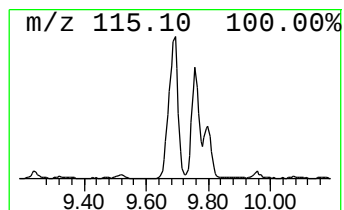
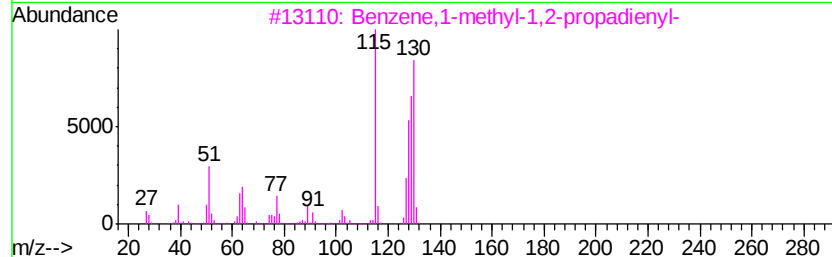
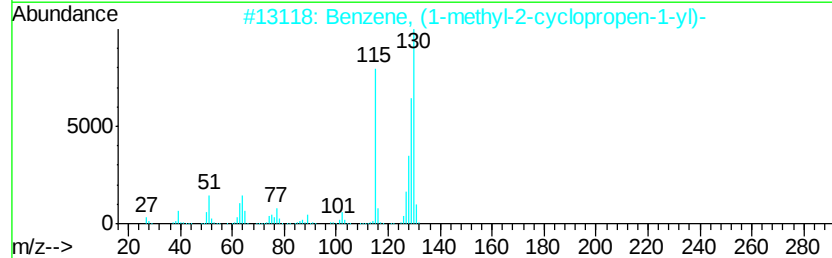
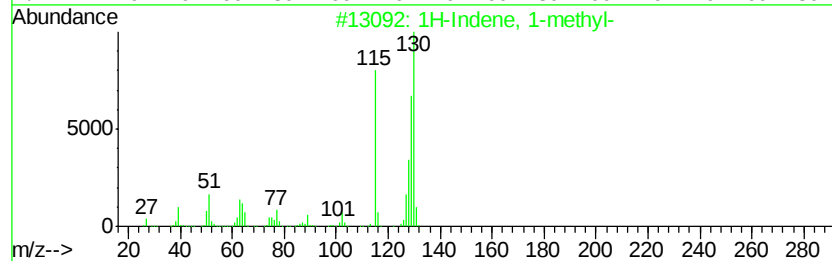
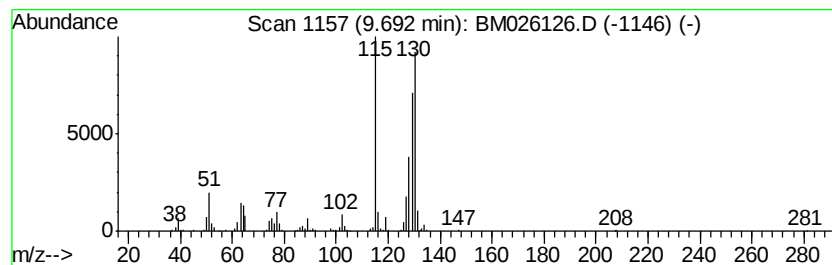
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 1H-Indene, 1-methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.69	7.05 ng/ul	508180	Naphthalene-d8	10.25

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1H-Indene, 1-methyl-	130	C10H10	000767-59-9	96
2		Benzene, (1-methyl-2-cyclopropen...	130	C10H10	065051-83-4	96
3		Benzene, 1-methyl-1,2-propadienyl-	130	C10H10	022433-39-2	96
4		2-Methylindene	130	C10H10	002177-47-1	96
5		1H-Indene, 3-methyl-	130	C10H10	000767-60-2	95



Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM052620\
Data File : BM026126.D
Acq On : 26 May 2020 23:05
Operator : MA/SJ
Sample : L2756-09
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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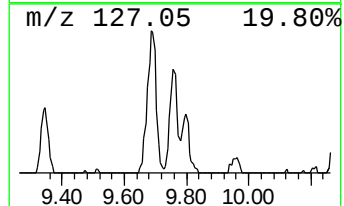
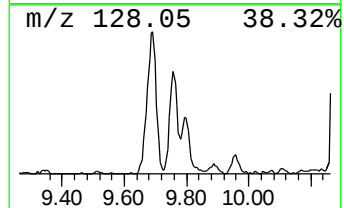
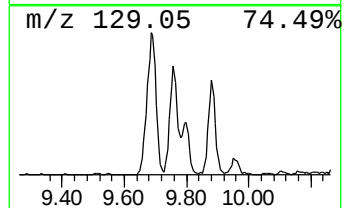
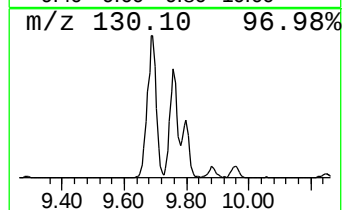
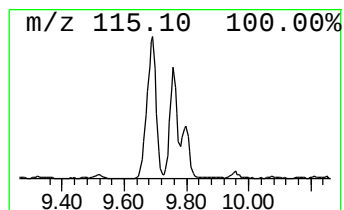
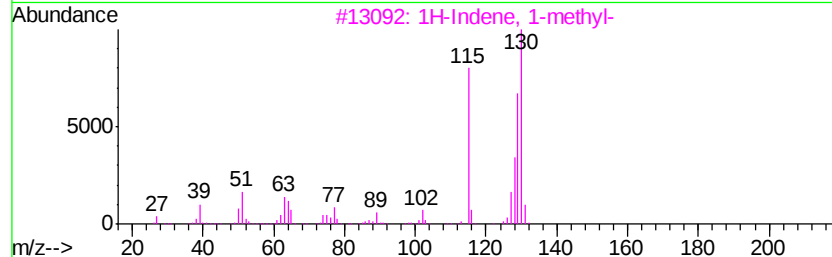
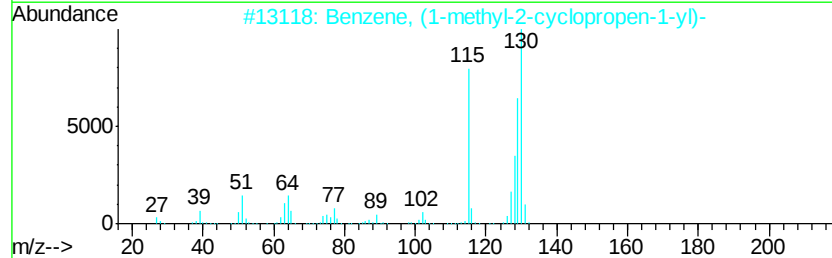
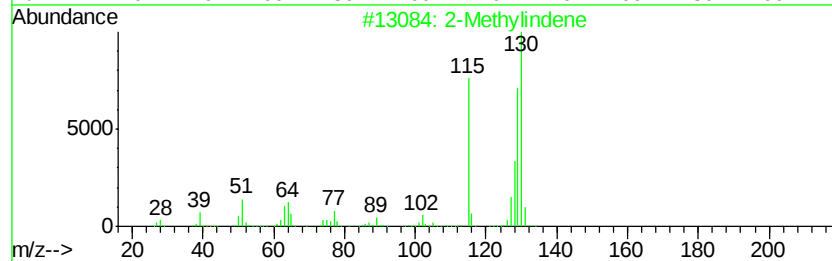
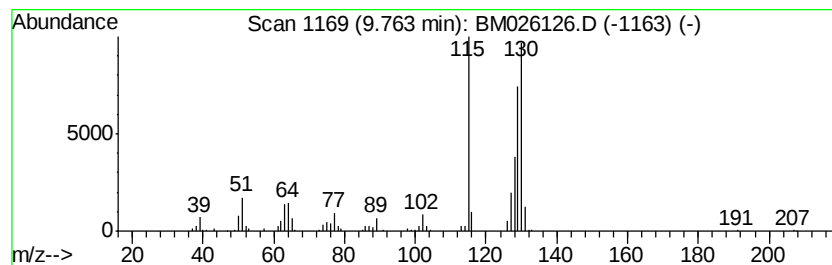
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 2-Methylindene Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.76	3.52 ng/ul	254057	Naphthalene-d8	10.25

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Methylindene	130	C10H10	002177-47-1	97
2			Benzene, (1-methyl-2-cyclopropen...	130	C10H10	065051-83-4	96
3			1H-Indene, 1-methyl-	130	C10H10	000767-59-9	96
4			Cycloprop[a]indene, 1,1a,6,6a-te...	130	C10H10	015677-15-3	95
5			1H-Indene, 3-methyl-	130	C10H10	000767-60-2	94



Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM052620\
Data File : BM026126.D
Acq On : 26 May 2020 23:05
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Sample : L2756-09
Misc :
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Instrument :
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ClientSampleId :
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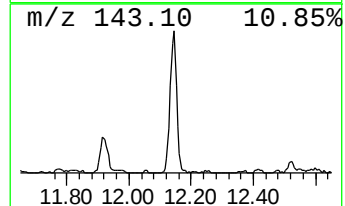
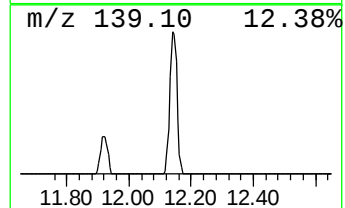
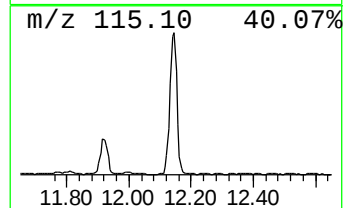
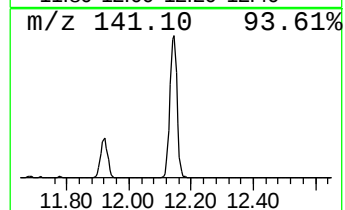
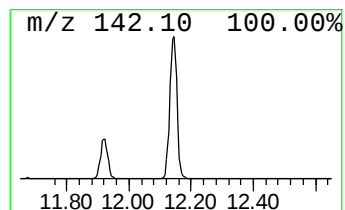
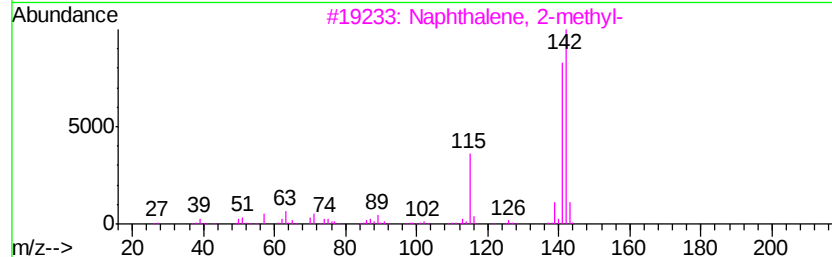
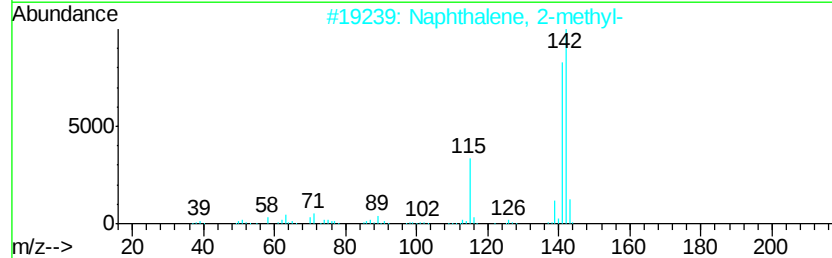
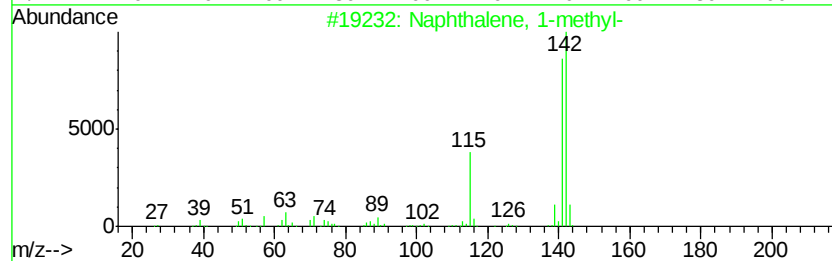
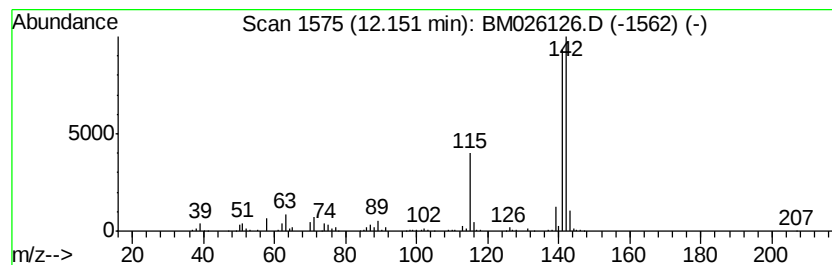
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Quant Title : SVOA CALIBRATION

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

Peak Number 11 Naphthalene, 1-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.15	9.64 ng/ul	694749	Naphthalene-d8	10.25

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 1-methyl-	142	C11H10	000090-12-0	96
2		Naphthalene, 2-methyl-	142	C11H10	000091-57-6	96
3		Naphthalene, 2-methyl-	142	C11H10	000091-57-6	96
4		1,4-Methanonaphthalene, 1,4-dihy...	142	C11H10	004453-90-1	94
5		Naphthalene, 1-methyl-	142	C11H10	000090-12-0	93



Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM052620\
Data File : BM026126.D
Acq On : 26 May 2020 23:05
Operator : MA/SJ
Sample : L2756-09
Misc :
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Instrument :
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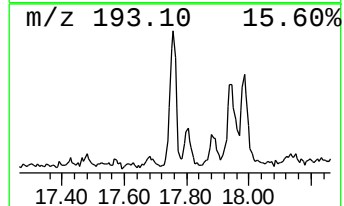
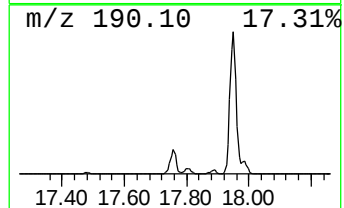
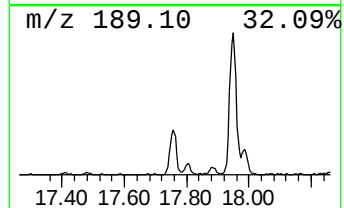
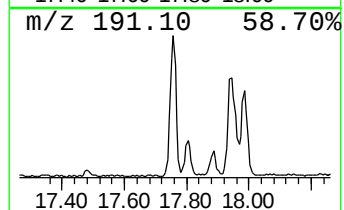
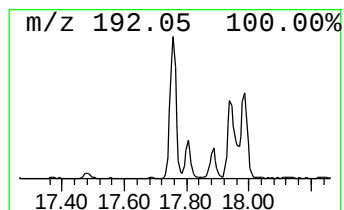
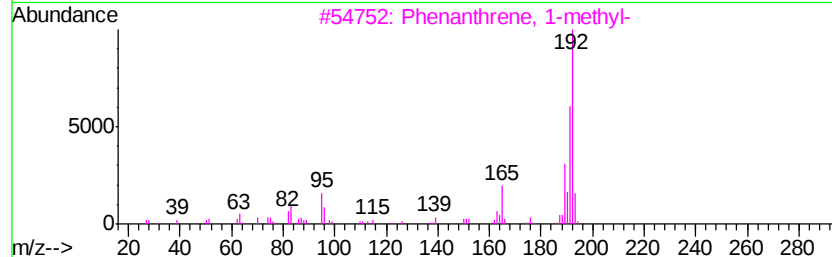
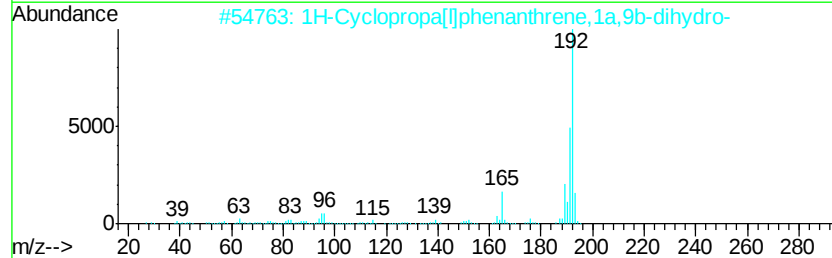
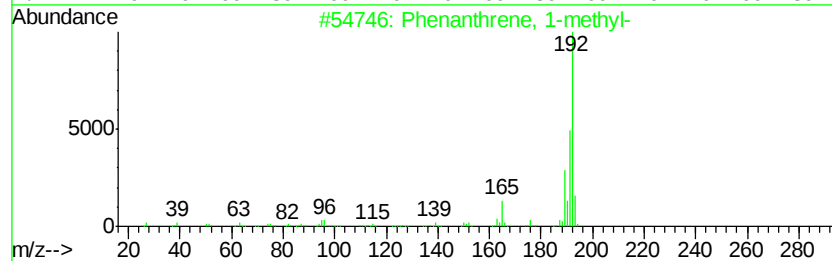
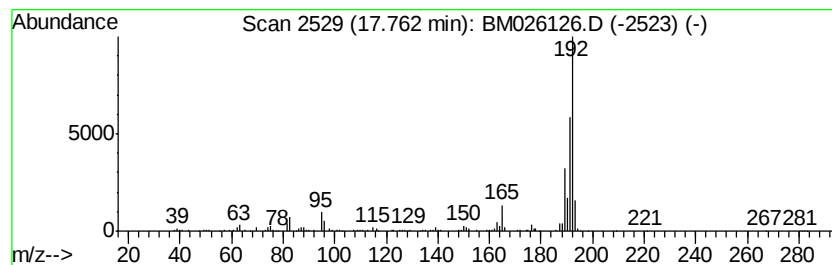
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Quant Title : SVOA CALIBRATION

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

Peak Number 12 Phenanthrene, 1-methyl- Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.76	2.26 ng/ul	237163	Phenanthrene-d10	16.88

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	96
2			1H-Cyclopropa[1]phenanthrene, 1a, ...	192	C15H12	000949-41-7	95
3			Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	94
4			Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	94
5			Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	93



Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM052620\
Data File : BM026126.D
Acq On : 26 May 2020 23:05
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Sample : L2756-09
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
ETKX1

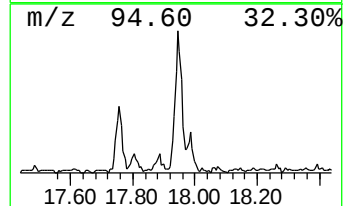
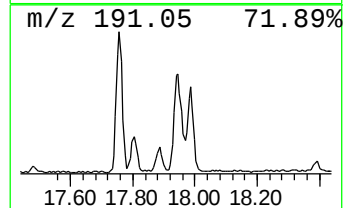
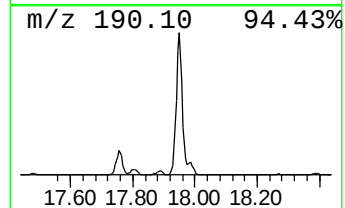
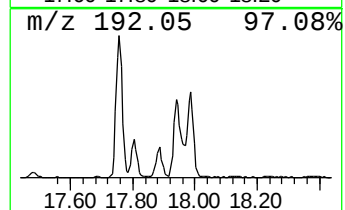
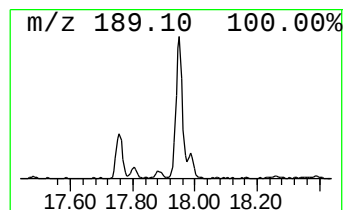
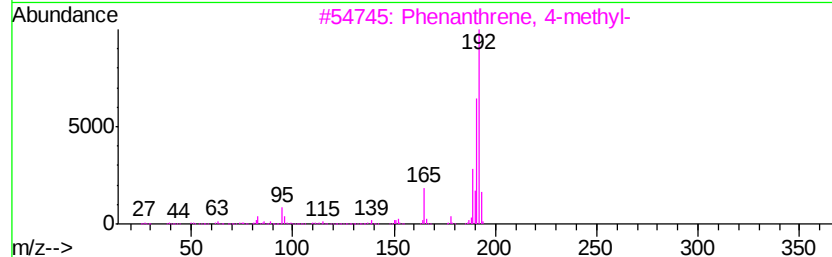
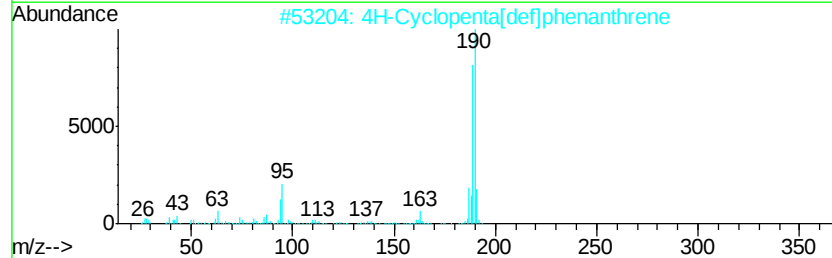
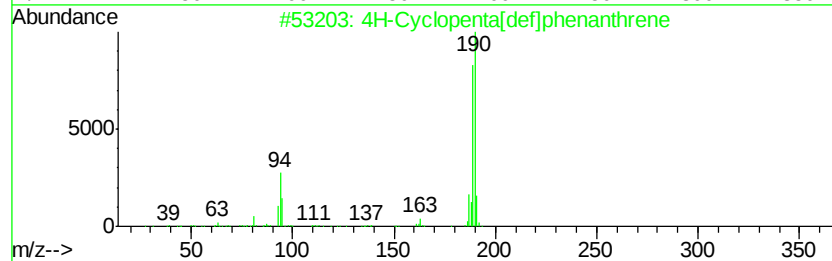
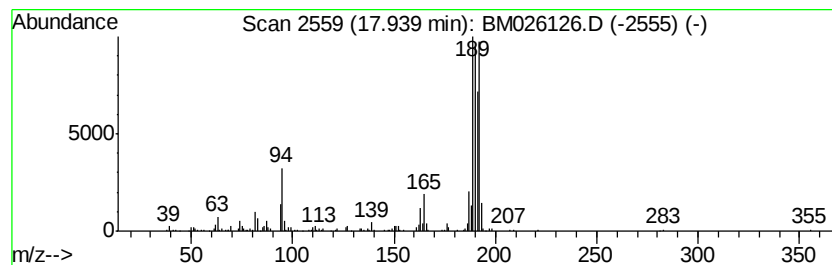
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Quant Title : SVOA CALIBRATION

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

Peak Number 13 4H-Cyclopenta[def]phenanthrene Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.94	3.55 ng/ul	373512	Phenanthrene-d10	16.88

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	58
2			4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	55
3			Phenanthrene, 4-methyl-	192	C15H12	000832-64-4	35
4			Phenanthrene, 4-methyl-	192	C15H12	000832-64-4	35
5			5H-Dibenzo[a,d]cycloheptene	192	C15H12	000256-81-5	18



Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM052620\
Data File : BM026126.D
Acq On : 26 May 2020 23:05
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Sample : L2756-09
Misc :
ALS Vial : 20 Sample Multiplier: 1

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BNA_M
ClientSampleId :
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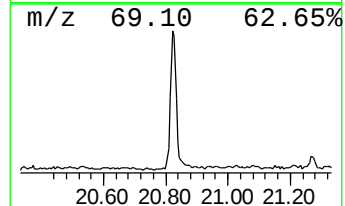
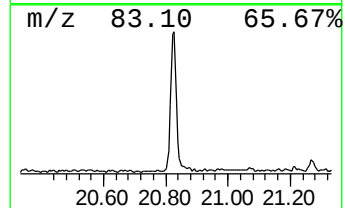
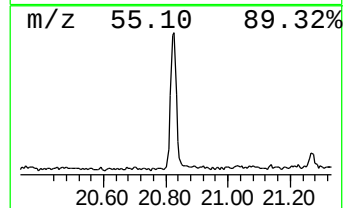
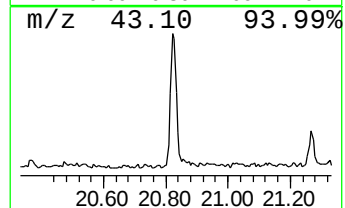
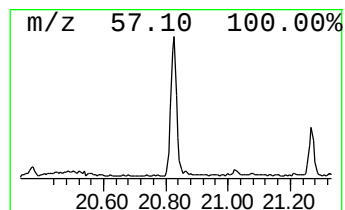
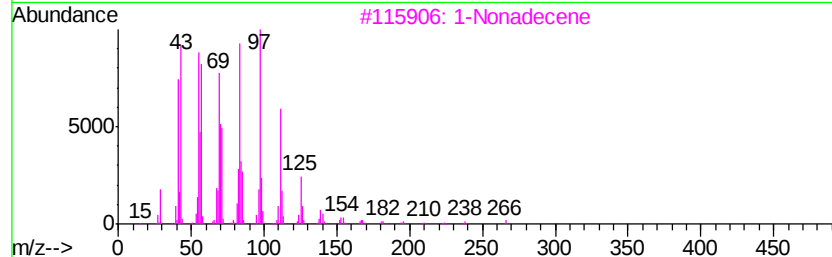
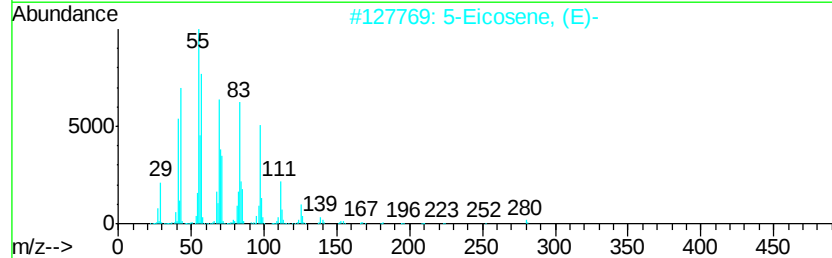
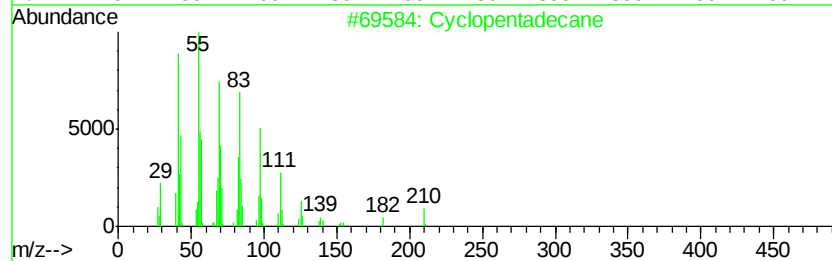
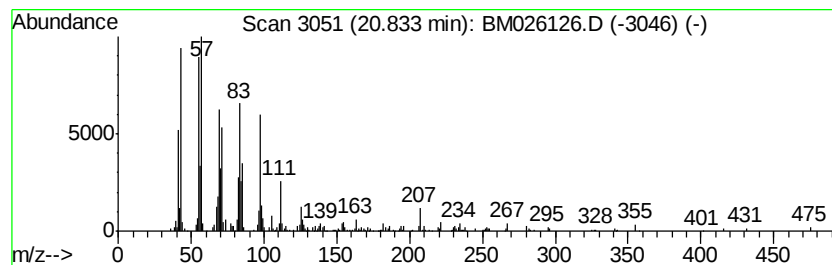
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Quant Title : SVOA CALIBRATION

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

Peak Number 14 (DEL) Alkane: Cyclic20.83 Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.83	2.30 ng/ul	285618	Chrysene-d12	21.08

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclopentadecane	210	C15H30	000295-48-7	96
2		5-Eicosene, (E)-	280	C20H40	074685-30-6	94
3		1-Nonadecene	266	C19H38	018435-45-5	94
4		Octacosanol	410	C28H58O	000557-61-9	94
5		1-Tricosene	322	C23H46	018835-32-0	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM052620\
Data File : BM026126.D
Acq On : 26 May 2020 23:05
Operator : MA/SJ
Sample : L2756-09
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
ETKX1

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM051320MA.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
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Benzene, 1-ethyl-...	7.60	3.5	ng/ul	170953	1	7.49	989909	20.0
Indene	8.02	42.6	ng/ul	2110860	1	7.49	989909	20.0
1H-Indene, 1-methyl-	9.69	7.0	ng/ul	508180	2	10.25	1441490	20.0
2-Methylindene	9.76	3.5	ng/ul	254057	2	10.25	1441490	20.0
Naphthalene, 1-me...	12.15	9.6	ng/ul	694749	2	10.25	1441490	20.0
Phenanthrene, 1-m...	17.76	2.3	ng/ul	237163	4	16.88	2102640	20.0
4H-Cyclopenta[def...	17.94	3.5	ng/ul	373512	4	16.88	2102640	20.0
(DEL) Alkane: Cyc...	20.83	2.3	ng/ul	285618	5	21.08	2479070	20.0