

Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM060120\
 Data File : BM026222.D
 Acq On : 02 Jun 2020 11:31
 Operator : JU/CG
 Sample : L2829-07DL 5X
 Misc :
 ALS Vial : 27 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C5CD2DL

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	5.516	442	447	453	rBV	48046	69616	0.15%	0.097%
2	6.669	639	643	652	rVB	79508	128260	0.28%	0.178%
3	6.822	662	669	674	rBV	103604	166549	0.37%	0.232%
4	6.875	674	678	687	rVB	44197	73568	0.16%	0.102%
5	7.010	696	701	709	rVB	126543	196723	0.44%	0.274%
6	7.128	716	721	727	rVB	174607	251337	0.56%	0.349%
7	7.469	772	779	785	rBV	674796	1014491	2.25%	1.410%
8	7.587	793	799	807	rVB	130538	201932	0.45%	0.281%
9	7.840	837	842	846	rBV	36160	54932	0.12%	0.076%
10	8.192	893	902	910	rVV	128844	217129	0.48%	0.302%
11	8.569	960	966	969	rBV	38699	57925	0.13%	0.081%
12	8.616	969	974	985	rVB2	123419	226028	0.50%	0.314%
13	9.087	1049	1054	1059	rVV	33274	48522	0.11%	0.067%
14	9.145	1059	1064	1074	rVB	40311	66429	0.15%	0.092%
15	9.328	1090	1095	1103	rVB	93157	159184	0.35%	0.221%
16	9.645	1136	1149	1158	rBV3	67731	171576	0.38%	0.239%
17	9.869	1180	1187	1195	rVB	171338	288260	0.64%	0.401%
18	10.139	1228	1233	1240	rBV4	36374	87083	0.19%	0.121%
19	10.228	1241	1248	1253	rVV	823610	1437926	3.19%	1.999%
20	10.281	1253	1257	1266	rVV	803307	1279627	2.84%	1.779%
21	10.386	1266	1275	1284	rVB3	72569	156964	0.35%	0.218%
22	10.651	1315	1320	1331	rVB3	15457	51592	0.11%	0.072%
23	11.904	1527	1533	1542	rVB	300263	512818	1.14%	0.713%
24	12.051	1549	1558	1562	rBV2	78710	158207	0.35%	0.220%
25	12.128	1565	1571	1576	rVV	280173	485558	1.08%	0.675%
26	12.169	1576	1578	1586	rVV6	29237	66843	0.15%	0.093%
27	12.245	1586	1591	1598	rVB	91246	160236	0.36%	0.223%
28	12.451	1621	1626	1633	rVB9	27700	51985	0.12%	0.072%
29	12.545	1639	1642	1646	rVV4	29782	53726	0.12%	0.075%
30	12.610	1647	1653	1657	rVV5	59704	121785	0.27%	0.169%
31	12.645	1657	1659	1665	rVB5	33166	48488	0.11%	0.067%
32	12.839	1684	1692	1702	rVV2	179179	356164	0.79%	0.495%
33	12.939	1706	1709	1715	rBV4	25514	51311	0.11%	0.071%
34	13.086	1729	1734	1738	rVB3	48115	71533	0.16%	0.099%

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35	13.139	1738	1743	1757	rVB3	72804	196775	0.44%	0.274%
36	13.251	1757	1762	1763	rBV2	55724	69356	0.15%	0.096%
37	13.351	1773	1779	1784	rVB4	64881	127452	0.28%	0.177%
38	13.433	1785	1793	1798	rBV3	116124	227587	0.50%	0.316%
39	13.492	1798	1803	1806	rBV2	76007	109743	0.24%	0.153%
40	13.539	1806	1811	1817	rVB	297132	430482	0.96%	0.599%
41	13.669	1829	1833	1837	rVV	66574	88342	0.20%	0.123%
42	13.710	1837	1840	1844	rVB4	41883	66264	0.15%	0.092%
43	13.804	1851	1856	1866	rVB	307851	554506	1.23%	0.771%
44	13.892	1867	1871	1874	rBV3	35541	54894	0.12%	0.076%
45	13.951	1877	1881	1893	rVB5	64639	140399	0.31%	0.195%
46	14.116	1903	1909	1923	rBV2	1166705	1853866	4.11%	2.578%
47	14.298	1936	1940	1943	rBV4	40419	61862	0.14%	0.086%
48	14.351	1946	1949	1955	rVB	61064	92595	0.21%	0.129%
49	14.404	1955	1958	1963	rVB6	42945	64307	0.14%	0.089%
50	14.527	1975	1979	1987	rVB5	49221	91217	0.20%	0.127%
51	14.674	1999	2004	2008	rVB	378128	533786	1.18%	0.742%
52	14.757	2008	2018	2023	rBV	1862498	2672564	5.93%	3.716%
53	14.798	2023	2025	2029	rVB3	95652	104021	0.23%	0.145%
54	15.016	2058	2062	2064	rBV3	31906	47276	0.10%	0.066%
55	15.069	2067	2071	2074	rVV3	53773	96490	0.21%	0.134%
56	15.116	2074	2079	2084	rVB	425724	595475	1.32%	0.828%
57	15.251	2098	2102	2104	rBV2	127716	178731	0.40%	0.248%
58	15.274	2104	2106	2119	rVB2	244118	408789	0.91%	0.568%
59	15.398	2123	2127	2129	rBV4	49945	72712	0.16%	0.101%
60	15.463	2135	2138	2147	rVB	162838	253131	0.56%	0.352%
61	15.574	2153	2157	2159	rBV5	38171	58472	0.13%	0.081%
62	15.874	2206	2208	2217	rVB5	55293	96648	0.21%	0.134%
63	15.998	2226	2229	2234	rVB5	32592	46150	0.10%	0.064%
64	16.157	2254	2256	2262	rVB2	37477	52907	0.12%	0.074%
65	16.551	2313	2323	2338	rBV	26554291	45068593	100.00%	62.661%
66	16.827	2366	2370	2372	rBV2	82228	97621	0.22%	0.136%
67	16.868	2372	2377	2382	rVV	1433668	1889381	4.19%	2.627%
68	16.910	2382	2384	2389	rVB3	64638	84178	0.19%	0.117%
69	16.968	2389	2394	2402	rBV	497311	771037	1.71%	1.072%
70	17.245	2439	2441	2447	rVB3	55430	71350	0.16%	0.099%
71	17.927	2553	2557	2558	rBV4	37177	50150	0.11%	0.070%

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72	18.092	2583	2585	2590	rVB5	49689	61647	0.14%	0.086%
73	19.268	2780	2785	2792	rVB	506581	689058	1.53%	0.958%
74	21.068	3086	3091	3098	rBV	1802473	2241918	4.97%	3.117%
75	23.044	3423	3427	3433	rVB	365587	576996	1.28%	0.802%
76	23.180	3444	3450	3461	rVB	1323817	2331615	5.17%	3.242%

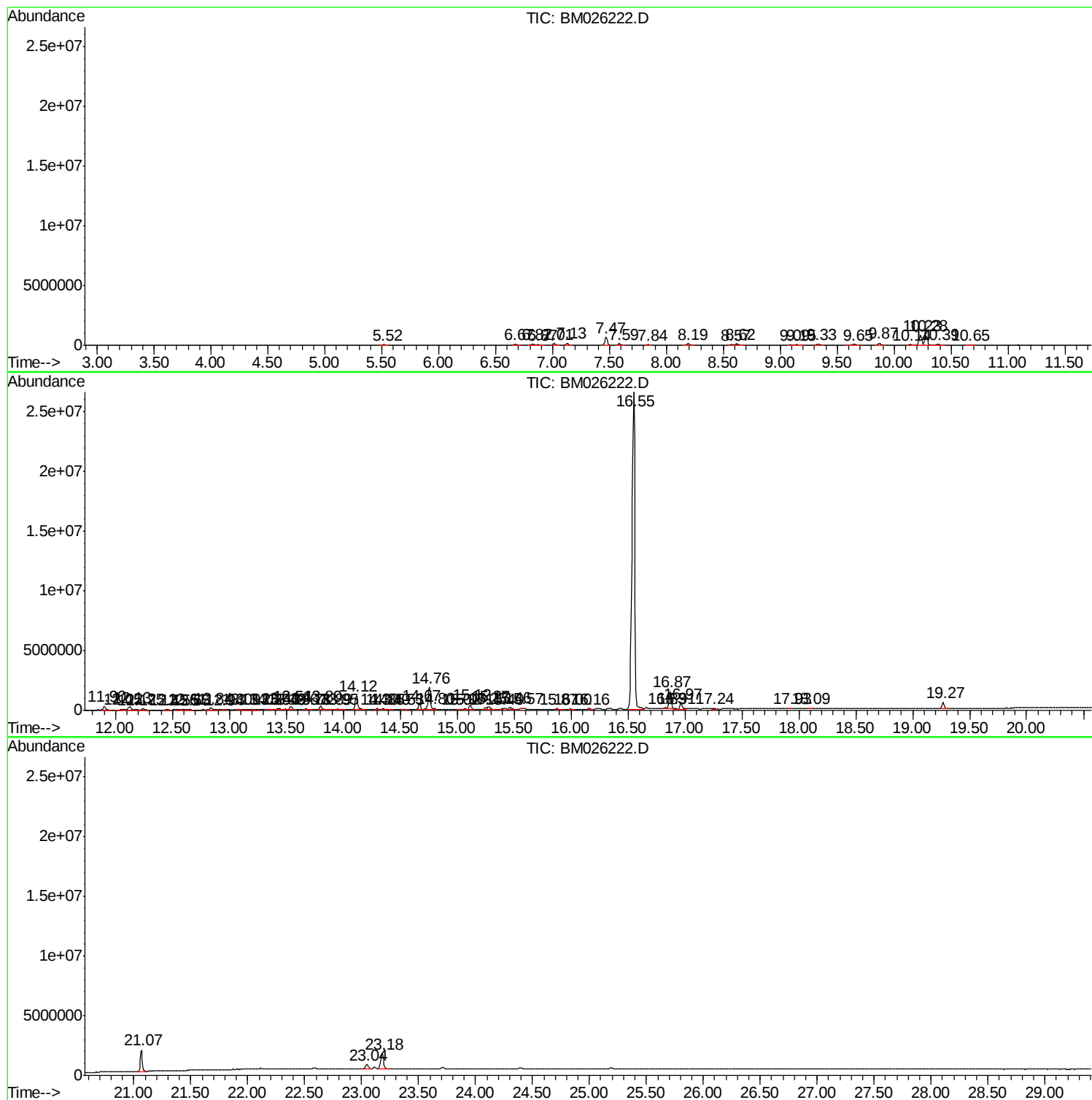
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Instrument :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM051320MA.M
Quant Title : SVOA CALIBRATION

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



Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM060120\
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Instrument :
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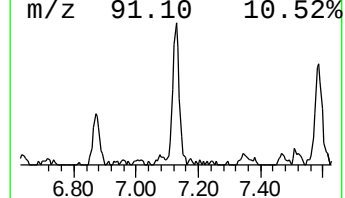
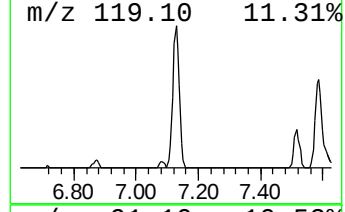
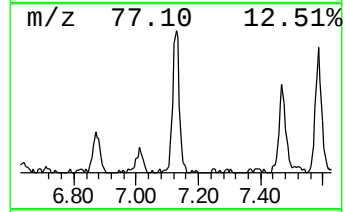
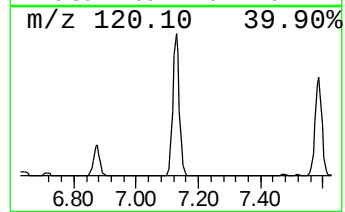
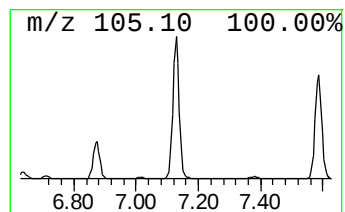
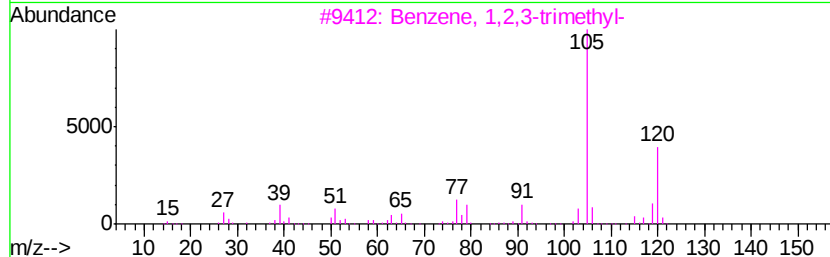
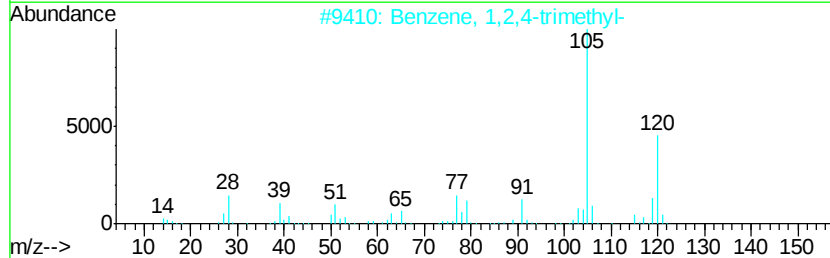
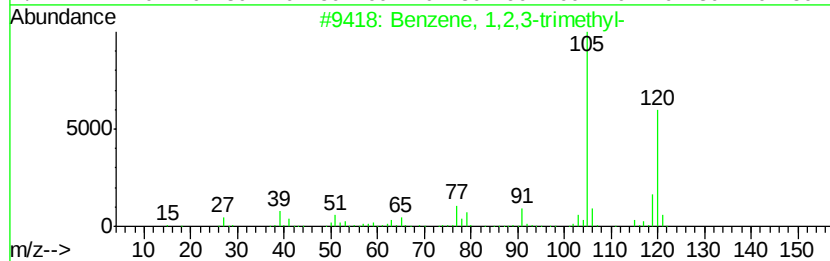
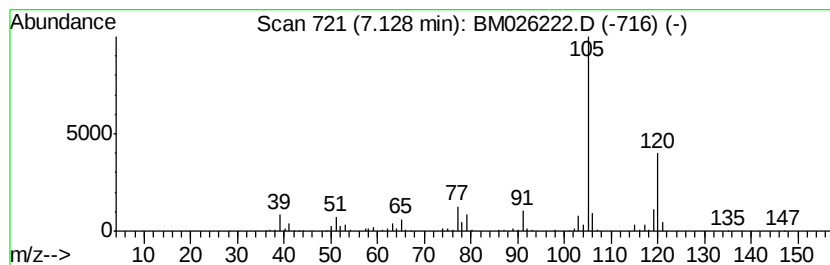
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM051320MA.M
Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.13	4.95 ng/ul	251337	1,4-Dichlorobenzene-d4	7.47

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	97
2			Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	94
3			Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	94
4			Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	93
5			Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91



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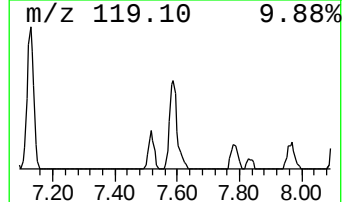
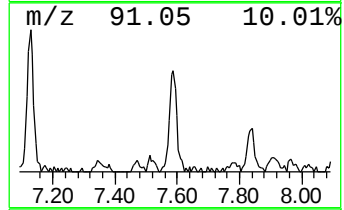
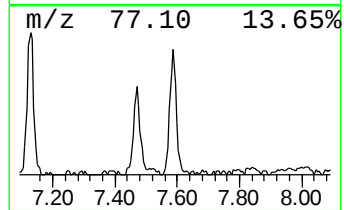
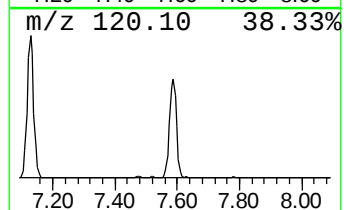
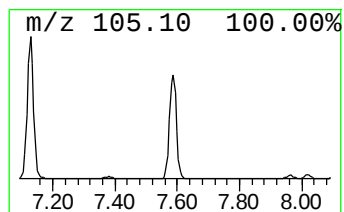
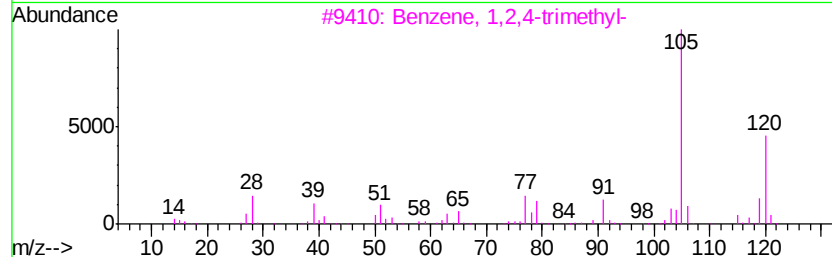
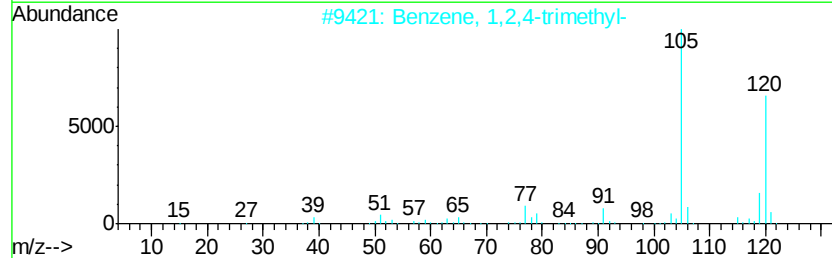
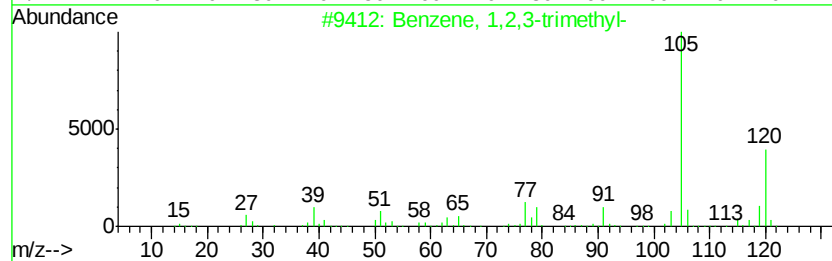
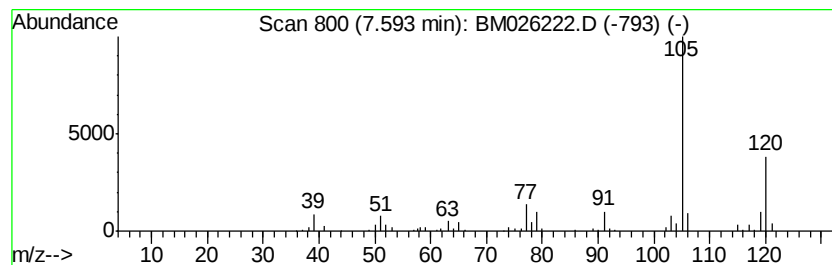
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM051320MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 Benzene, 1,2,4-trimethyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.59	3.98 ng/ul	201932	1,4-Dichlorobenzene-d4	7.47

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	97
2		Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	94
3		Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91
4		Mesitylene	120	C9H12	000108-67-8	91
5		Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91



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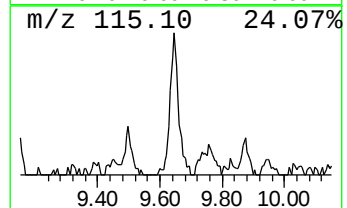
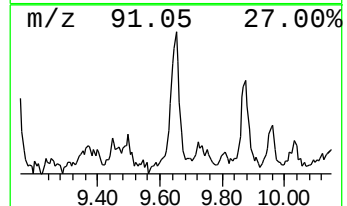
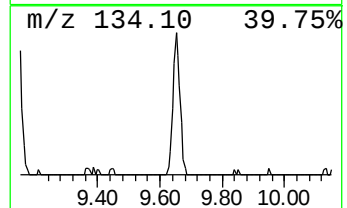
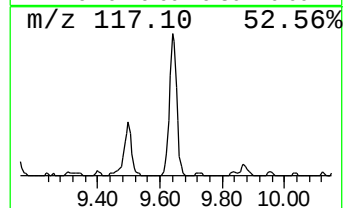
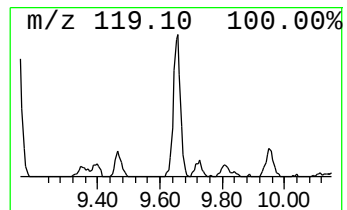
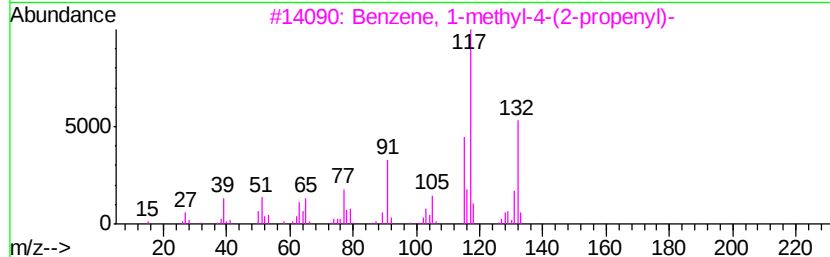
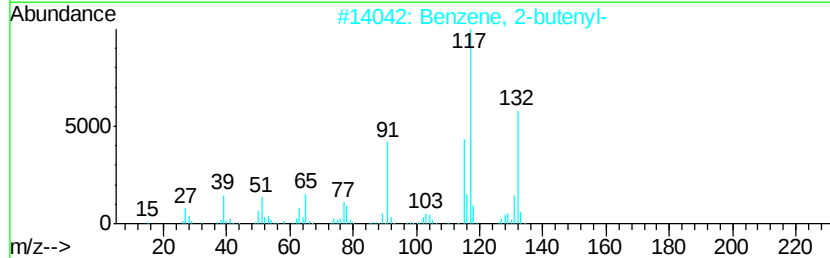
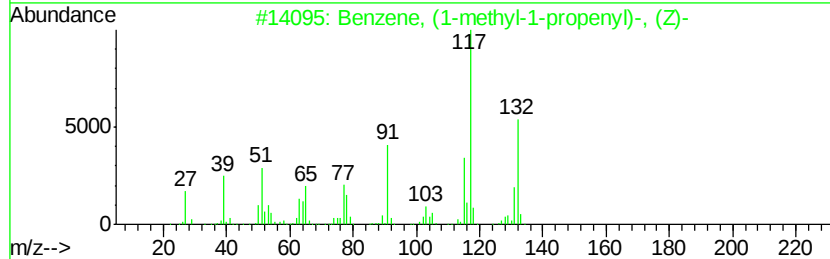
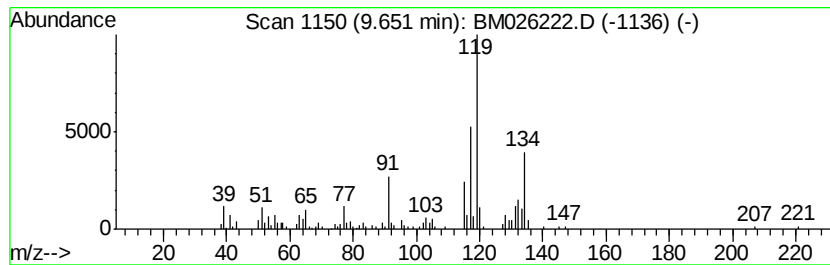
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 3 Benzene, (1-methyl-1-propen... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.65	2.39 ng/ul	171576	Naphthalene-d8	10.23

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene,	(1-methyl-1-propenyl)-,...	132	C10H12	000767-99-7	70	
2	Benzene,	2-butenyl-	132	C10H12	001560-06-1	70	
3	Benzene,	1-methyl-4-(2-propenyl)-	132	C10H12	003333-13-9	70	
4	Benzene,	(1-methyl-1-propenyl)-,...	132	C10H12	000768-00-3	55	
5	1-Phenyl-1-butene		132	C10H12	000824-90-8	55	



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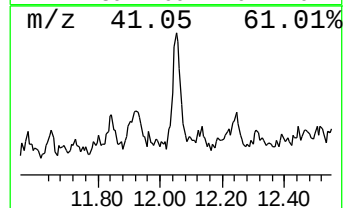
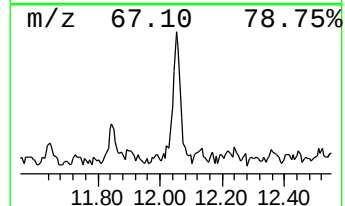
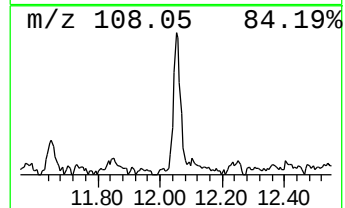
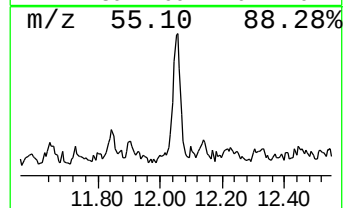
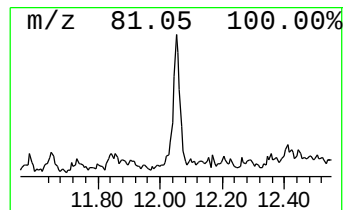
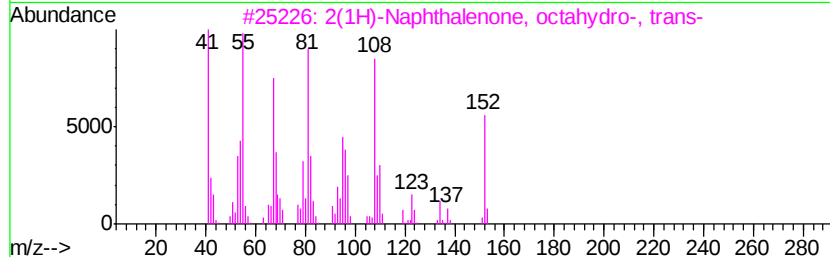
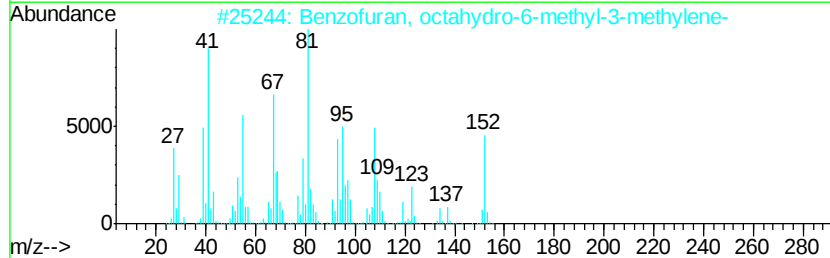
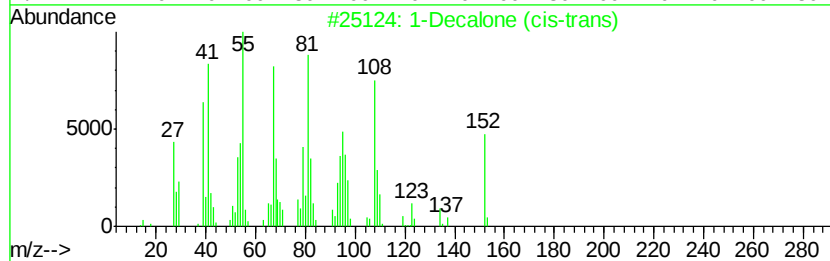
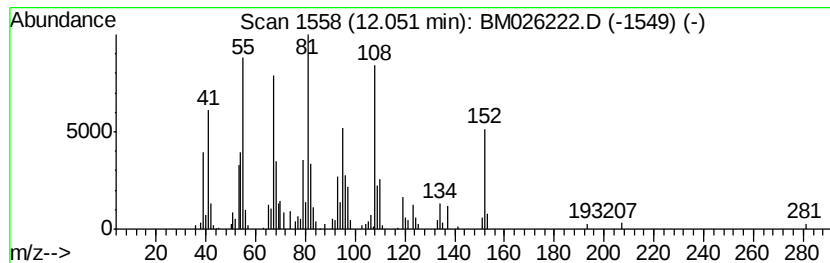
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ClientSampleId :
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 1-Decalone (cis-trans) Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.		
12.05	2.20 ng/ul	158207	Naphthalene-d8	10.23		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Decalone (cis-trans)		152	C10H16O	004832-16-0	93
2	Benzofuran, octahydro-6-methyl-3...		152	C10H16O	077954-13-3	90
3	2(1H)-Naphthalenone, octahydro-, ...		152	C10H16O	016021-08-2	72
4	Spiro[3.5]nonan-1-one, 5-methyl-...		152	C10H16O	065147-56-0	68
5	2-Decalone, c&t		152	C10H16O	004832-17-1	68



Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM060120\
Data File : BM026222.D
Acq On : 02 Jun 2020 11:31
Operator : JU/CG
Sample : L2829-07DL 5X
Misc :
ALS Vial : 27 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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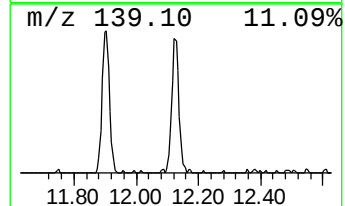
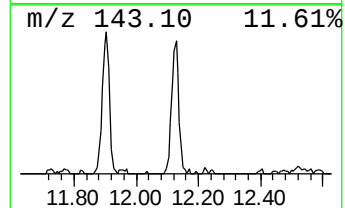
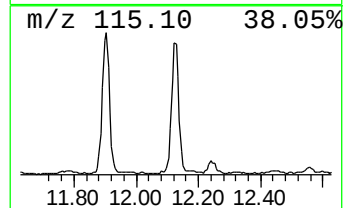
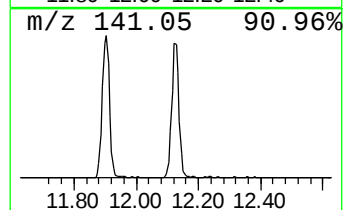
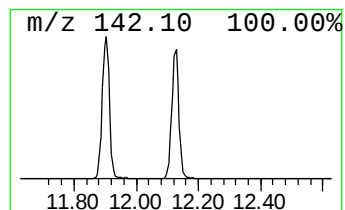
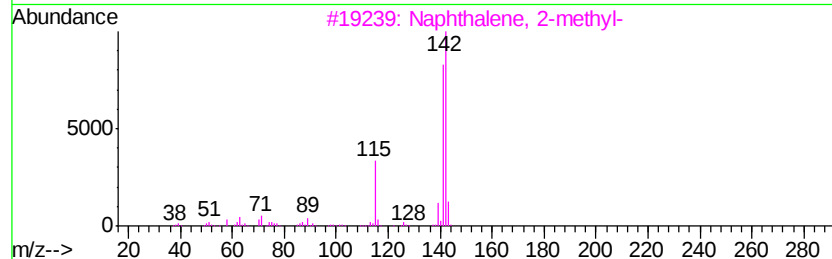
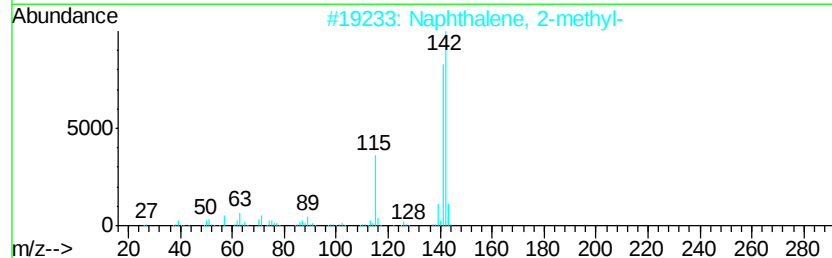
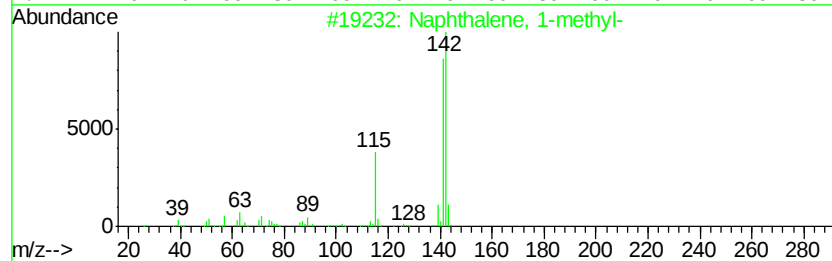
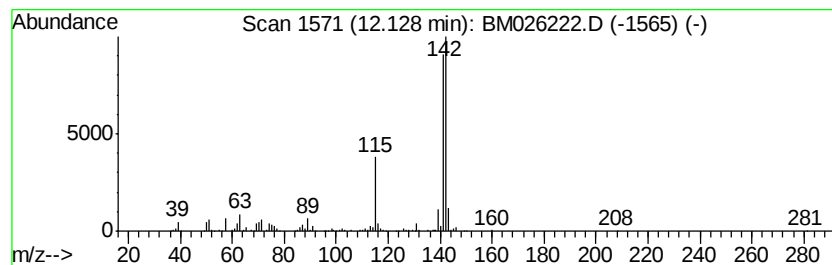
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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.13	6.75 ng/ul	485558	Naphthalene-d8	10.23

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	95
4		1,4-Methanonaphthalene, 1,4-dihy...	142	C11H10	004453-90-1	94
5		Naphthalene, 1-methyl-	142	C11H10	000090-12-0	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM060120\
Data File : BM026222.D
Acq On : 02 Jun 2020 11:31
Operator : JU/CG
Sample : L2829-07DL 5X
Misc :
ALS Vial : 27 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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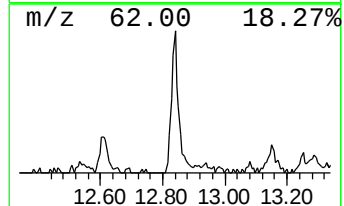
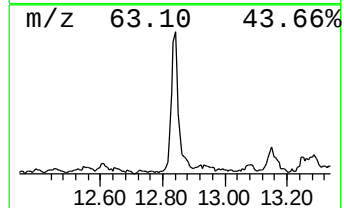
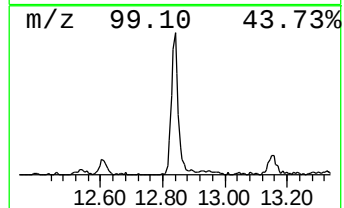
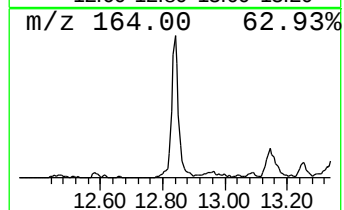
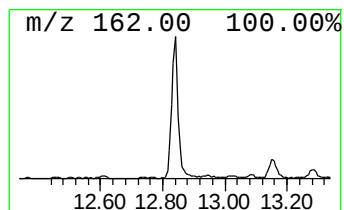
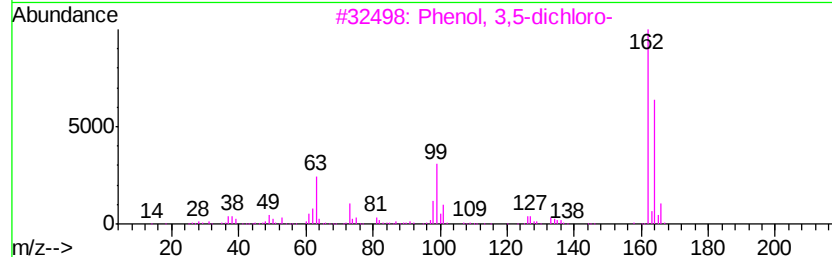
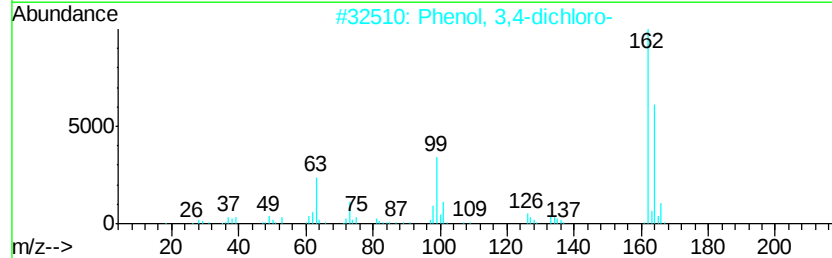
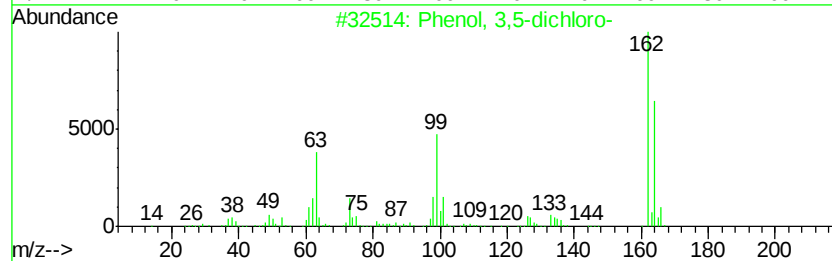
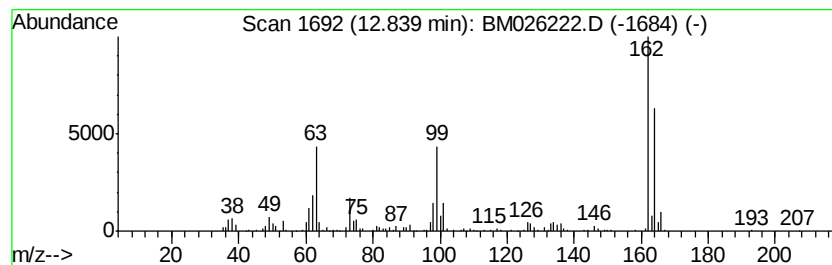
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 Phenol, 3,5-dichloro- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD		R.T.
12.84	3.84 ng/ul	356164	Acenaphthene-d10		14.12

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol,	3,5-dichloro-		162	C6H4Cl2O	000591-35-5	98
2	Phenol,	3,4-dichloro-		162	C6H4Cl2O	000095-77-2	97
3	Phenol,	3,5-dichloro-		162	C6H4Cl2O	000591-35-5	97
4	Phenol,	3,4-dichloro-		162	C6H4Cl2O	000095-77-2	97
5	Phenol,	3,4-dichloro-		162	C6H4Cl2O	000095-77-2	92



Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM060120\
Data File : BM026222.D
Acq On : 02 Jun 2020 11:31
Operator : JU/CG
Sample : L2829-07DL 5X
Misc :
ALS Vial : 27 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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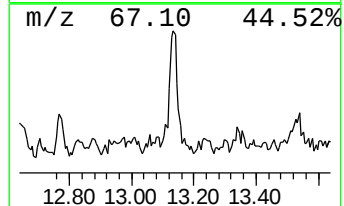
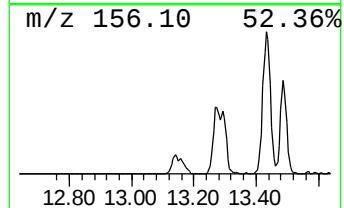
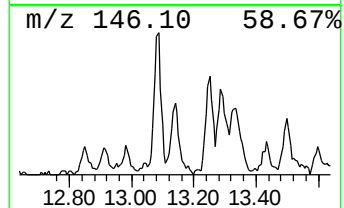
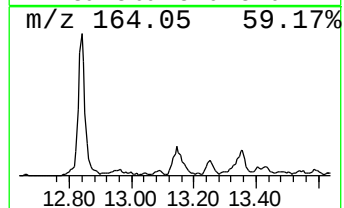
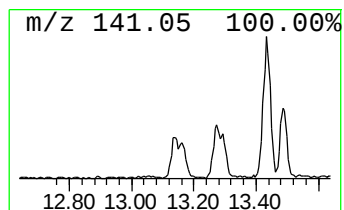
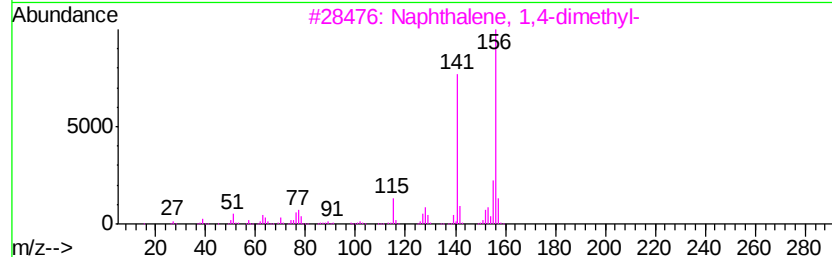
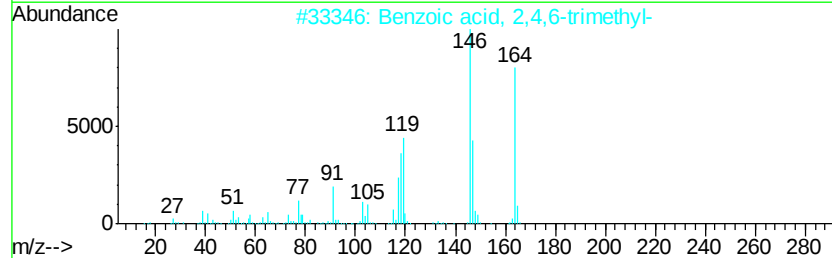
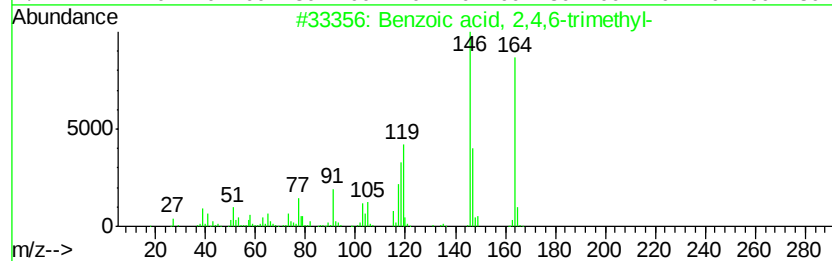
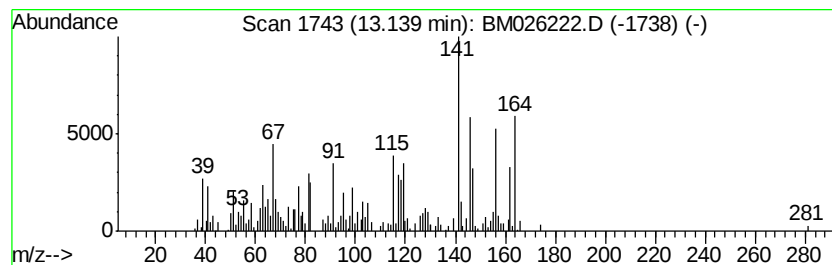
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 Benzoic acid, 2,4,6-trimethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.14	2.12 ng/ul	196775	Acenaphthene-d10	14.12

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzoic acid, 2,4,6-trimethyl-	164	C10H12O2	000480-63-7	66
2			Benzoic acid, 2,4,6-trimethyl-	164	C10H12O2	000480-63-7	60
3			Naphthalene, 1,4-dimethyl-	156	C12H12	000571-58-4	38
4			Naphthalene, 1,2-dimethyl-	156	C12H12	000573-98-8	38
5			Naphthalene, 1,3-dimethyl-	156	C12H12	000575-41-7	38



Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM060120\
Data File : BM026222.D
Acq On : 02 Jun 2020 11:31
Operator : JU/CG
Sample : L2829-07DL 5X
Misc :
ALS Vial : 27 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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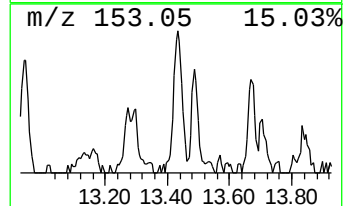
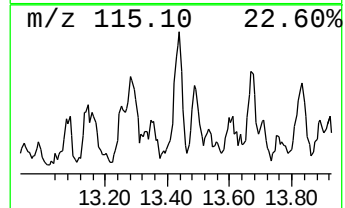
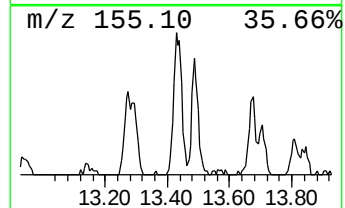
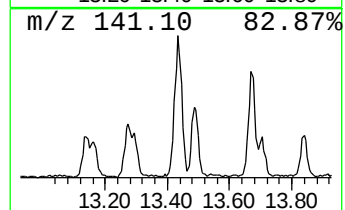
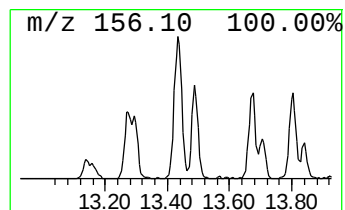
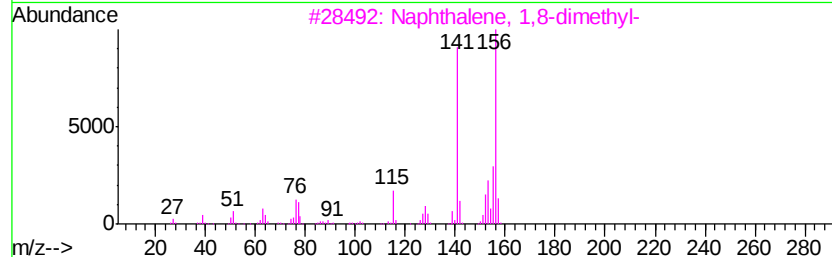
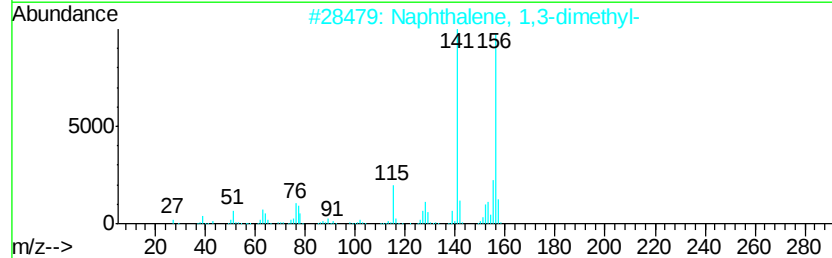
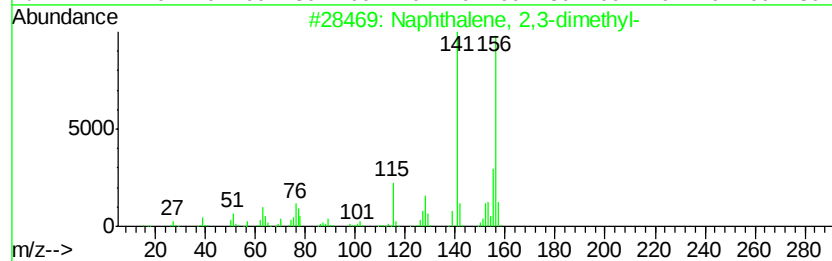
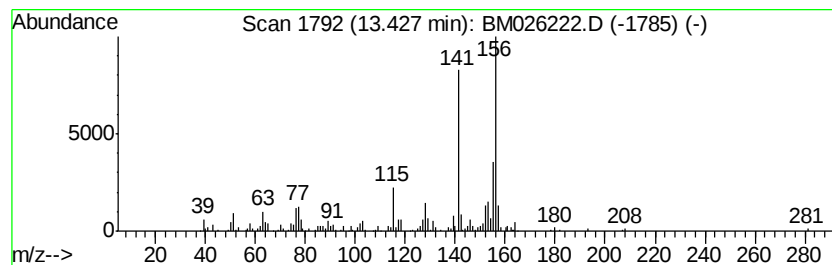
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 Naphthalene, 2,3-dimethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.43	2.46 ng/ul	227587	Acenaphthene-d10	14.12

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	98
2		Naphthalene, 1,3-dimethyl-	156	C12H12	000575-41-7	98
3		Naphthalene, 1,8-dimethyl-	156	C12H12	000569-41-5	98
4		Naphthalene, 1,7-dimethyl-	156	C12H12	000575-37-1	98
5		Naphthalene, 1,2-dimethyl-	156	C12H12	000573-98-8	98



Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM060120\
Data File : BM026222.D
Acq On : 02 Jun 2020 11:31
Operator : JU/CG
Sample : L2829-07DL 5X
Misc :
ALS Vial : 27 Sample Multiplier: 1

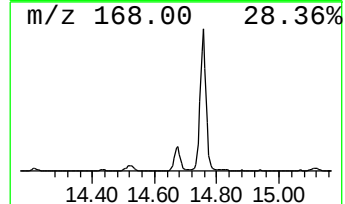
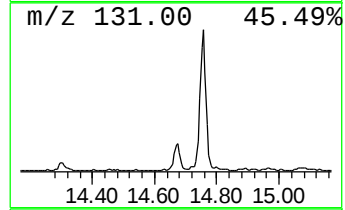
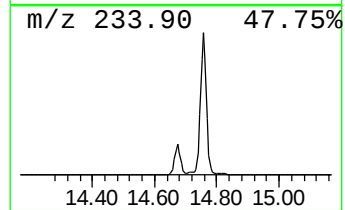
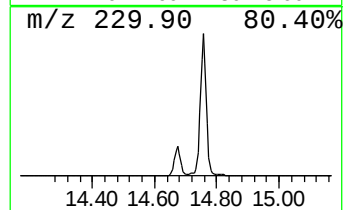
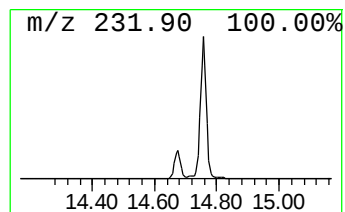
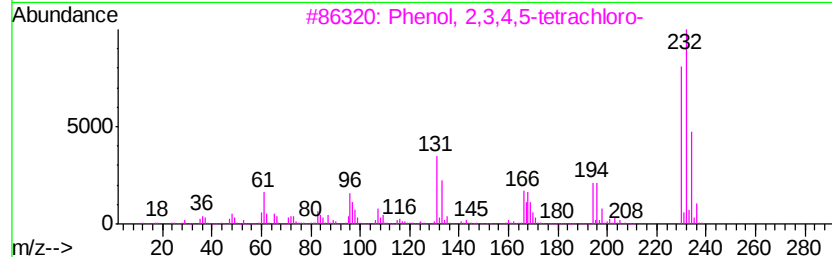
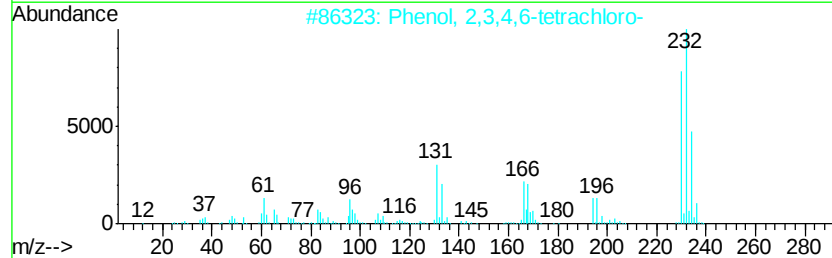
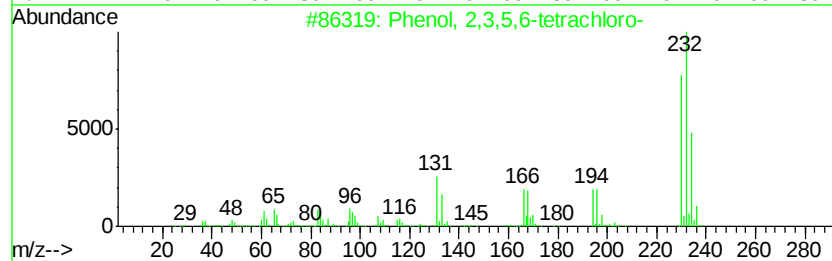
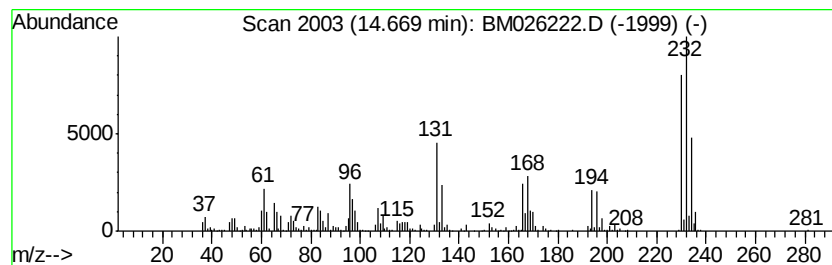
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ClientSampleId :
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 Phenol, 2,3,5,6-tetrachloro- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD		R.T.	
14.67	5.76 ng/ul	533786	Acenaphthene-d10		14.12	
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol,	2,3,5,6-tetrachloro-	230	C6H2Cl4O	000935-95-5	99
2	Phenol,	2,3,4,6-tetrachloro-	230	C6H2Cl4O	000058-90-2	99
3	Phenol,	2,3,4,5-tetrachloro-	230	C6H2Cl4O	004901-51-3	99
4	Phenol,	2,3,4,6-tetrachloro-	230	C6H2Cl4O	000058-90-2	98
5	Phenol,	2,3,5,6-tetrachloro-	230	C6H2Cl4O	000935-95-5	93



Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM060120\
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Misc :
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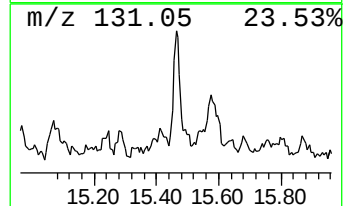
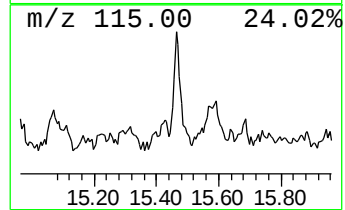
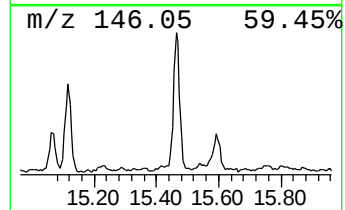
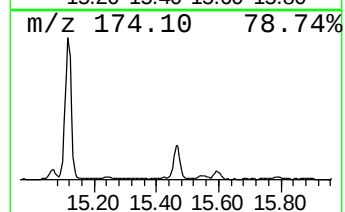
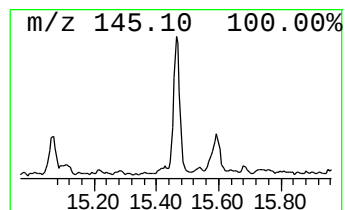
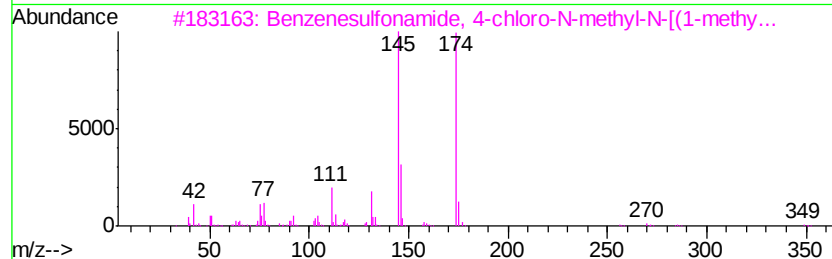
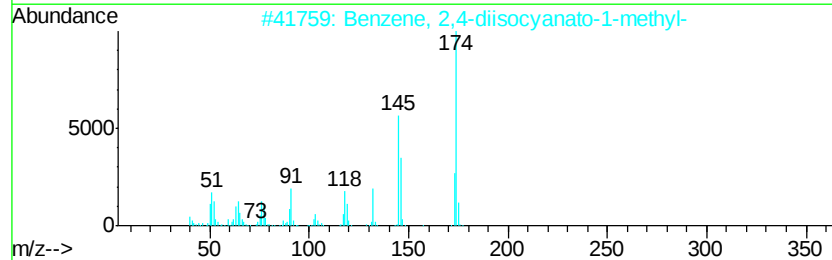
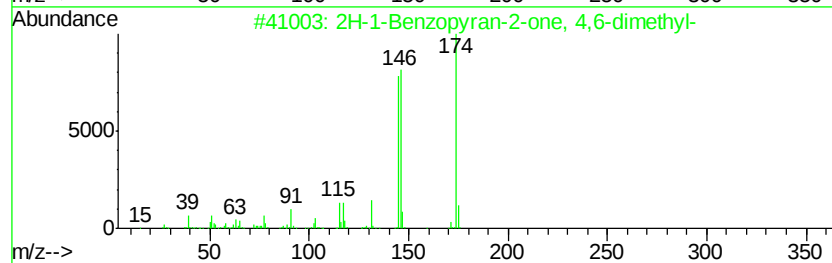
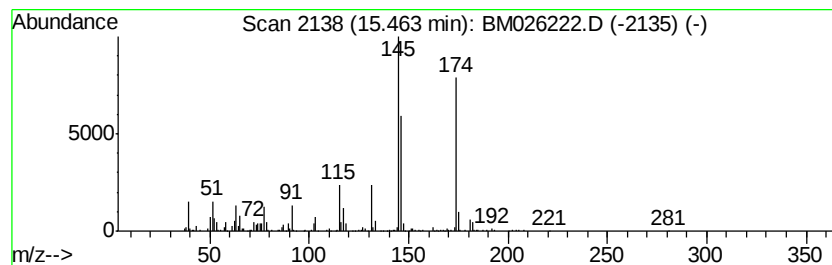
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 2H-1-Benzopyran-2-one, 4,6-... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.46	2.73 ng/ul	253131	Acenaphthene-d10	14.12

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2H-1-Benzopyran-2-one, 4,6-dimet...	174	C11H10O2	014002-89-2	70
2			Benzene, 2,4-diisocyanato-1-methyl-	174	C9H6N2O2	000584-84-9	64
3			Benzenesulfonamide, 4-chloro-N-m...	349	C16H16ClN3O2S	1000319-71-4	59
4			1,2-Dihydro-1,3-dimethyl-2-oxoqu...	174	C10H10N2O	003149-25-5	55
5			4-Dimethylaminopyrido[3,2-c]pyri...	174	C9H10N4	1000326-90-5	50



Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM060120\
 Data File : BM026222.D
 Acq On : 02 Jun 2020 11:31
 Operator : JU/CG
 Sample : L2829-07DL 5X
 Misc :
 ALS Vial : 27 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C5CD2DL

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
Benzene, 1,2,3-tr...	7.13	5.0	ng/ul	251337	1	7.47	1014490	20.0
Benzene, 1,2,4-tr...	7.59	4.0	ng/ul	201932	1	7.47	1014490	20.0
Benzene, (1-methy...	9.65	2.4	ng/ul	171576	2	10.23	1437930	20.0
1-Decalone (cis-t...	12.05	2.2	ng/ul	158207	2	10.23	1437930	20.0
Naphthalene, 1-me...	12.13	6.8	ng/ul	485558	2	10.23	1437930	20.0
Phenol, 3,5-dichl...	12.84	3.8	ng/ul	356164	3	14.12	1853870	20.0
Benzoic acid, 2,4...	13.14	2.1	ng/ul	196775	3	14.12	1853870	20.0
Naphthalene, 2,3-...	13.43	2.5	ng/ul	227587	3	14.12	1853870	20.0
Phenol, 2,3,5,6-t...	14.67	5.8	ng/ul	533786	3	14.12	1853870	20.0
2H-1-Benzopyran-2...	15.46	2.7	ng/ul	253131	3	14.12	1853870	20.0