

Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN062918\
 Data File : BN001921.D
 Acq On : 29 Jun 2018 09:58
 Operator : JU/SJ
 Sample : J3661-02
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 X2D20

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_N\METHODS\SOM-EPA-BN061918.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	3.023	3	6	28	rVB	1539337	2425358	4.47%	0.327%
2	3.270	44	48	60	rVB	128299	202535	0.37%	0.027%
3	3.805	135	139	146	rVB	49558	78042	0.14%	0.011%
4	3.981	164	169	178	rBV	305877	490662	0.90%	0.066%
5	4.875	314	321	328	rBV	148500	232131	0.43%	0.031%
6	5.869	486	490	498	rVB	41362	67109	0.12%	0.009%
7	6.911	655	667	679	rBV2	5141644	12126438	22.33%	1.635%
8	7.052	684	691	702	rVV	2208617	3367992	6.20%	0.454%
9	7.146	702	707	714	rVV	46405	81906	0.15%	0.011%
10	7.246	716	724	726	rVV	2672713	4084287	7.52%	0.551%
11	7.281	726	730	738	rVV	4896399	7992736	14.72%	1.078%
12	7.346	738	741	749	rVB4	36482	64991	0.12%	0.009%
13	7.599	775	784	794	rVB	4128698	6584212	12.13%	0.888%
14	7.711	796	803	805	rBV	1879628	2970388	5.47%	0.401%
15	7.746	805	809	820	rVB	5476847	8712477	16.05%	1.175%
16	8.058	851	862	875	rVB	5505709	9061427	16.69%	1.222%
17	8.240	883	893	898	rBV2	1477950	3715021	6.84%	0.501%
18	8.293	898	902	910	rVB	877255	1420424	2.62%	0.192%
19	8.411	910	922	938	rBV	3470299	5539903	10.20%	0.747%
20	8.781	977	985	991	rBV	4298431	7202861	13.26%	0.971%
21	8.858	991	998	1006	rVV	2657523	4309260	7.94%	0.581%
22	9.569	1110	1119	1130	rBV	2075065	3504387	6.45%	0.473%
23	9.681	1133	1138	1141	rVV2	39446	68247	0.13%	0.009%
24	9.716	1141	1144	1150	rVB	63772	93494	0.17%	0.013%
25	10.134	1203	1215	1238	rBV2	6751033	16324091	30.06%	2.201%
26	10.352	1244	1252	1266	rVB	5956306	10007653	18.43%	1.349%
27	10.487	1266	1275	1278	rBV	2843603	4826760	8.89%	0.651%
28	10.534	1278	1283	1291	rVV	8250268	14069763	25.91%	1.897%
29	10.616	1291	1297	1314	rVB	3298635	5769758	10.63%	0.778%
30	10.822	1325	1332	1347	rVB	5631211	9151258	16.85%	1.234%
31	11.199	1390	1396	1403	rVB	56001	84787	0.16%	0.011%
32	11.528	1443	1452	1467	rBV	5293979	9592113	17.66%	1.293%
33	12.016	1527	1535	1540	rBV	128764	204018	0.38%	0.028%
34	12.075	1540	1545	1552	rVB	70277	114129	0.21%	0.015%

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35	12.522	1613	1621	1629	rBV	10718454	17108979	31.51%	2.307%
36	12.751	1646	1660	1664	rBV	2779189	5653835	10.41%	0.762%
37	12.799	1664	1668	1682	rVB	2377308	4419723	8.14%	0.596%
38	13.757	1825	1831	1842	rVB	6352202	9040066	16.65%	1.219%
39	13.922	1851	1859	1870	rBV	7662169	10862308	20.00%	1.465%
40	14.028	1870	1877	1887	rVB	7001990	10958229	20.18%	1.478%
41	14.340	1923	1930	1935	rBV2	4247029	6687799	12.32%	0.902%
42	14.404	1935	1941	1949	rVB	14278853	21551078	39.69%	2.906%
43	14.545	1958	1965	1981	rBV	3220022	4979090	9.17%	0.671%
44	14.710	1986	1993	2001	rVV	7125814	9960047	18.34%	1.343%
45	15.175	2065	2072	2081	rBV	13166065	17929794	33.02%	2.418%
46	15.340	2092	2100	2104	rBV	8723038	13540908	24.94%	1.826%
47	15.393	2104	2109	2115	rVV	14985953	21778799	40.11%	2.937%
48	15.451	2115	2119	2131	rVB	3079421	4269569	7.86%	0.576%
49	15.575	2136	2140	2145	rVV	119879	178914	0.33%	0.024%
50	15.640	2146	2151	2157	rVB	334263	475673	0.88%	0.064%
51	16.251	2250	2255	2259	rBV	142268	185745	0.34%	0.025%
52	16.298	2259	2263	2270	rVV2	54023	83057	0.15%	0.011%
53	16.392	2272	2279	2287	rBV	11972611	17639574	32.49%	2.379%
54	16.739	2331	2338	2349	rBV	9923201	14175280	26.11%	1.911%
55	16.869	2358	2360	2367	rVB	40956	57753	0.11%	0.008%
56	17.087	2390	2397	2407	rVB	5623075	8288387	15.26%	1.118%
57	17.187	2407	2414	2417	rBV2	10373027	16363105	30.13%	2.206%
58	17.222	2417	2420	2431	rVB	15847674	22197928	40.88%	2.993%
59	17.704	2497	2502	2507	rBV	215228	303565	0.56%	0.041%
60	18.498	2634	2637	2641	rBV4	51576	76318	0.14%	0.010%
61	19.116	2737	2742	2751	rVB	141216	232183	0.43%	0.031%
62	19.392	2780	2789	2794	rBV2	310543	582350	1.07%	0.079%
63	19.486	2798	2805	2807	rVV2	12468924	18246633	33.60%	2.460%
64	19.516	2807	2810	2820	rVB	21650922	31280818	57.61%	4.218%
65	19.633	2826	2830	2835	rVB2	63526	84632	0.16%	0.011%
66	20.139	2912	2916	2922	rVV	314464	381527	0.70%	0.051%
67	20.198	2922	2926	2930	rVB	109040	135839	0.25%	0.018%
68	20.433	2961	2966	2971	rBV	20412124	24984308	46.01%	3.369%
69	20.586	2986	2992	2997	rBV	22834292	28490318	52.47%	3.842%
70	20.639	2998	3001	3005	rVB3	110656	142454	0.26%	0.019%
71	20.927	3047	3050	3054	rBV	106251	139577	0.26%	0.019%

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72	21.010	3061	3064	3068	rVB3	47452	59642	0.11%	0.008%
73	21.222	3095	3100	3104	rBV	27517866	29857291	54.99%	4.026%
74	21.275	3104	3109	3114	rVV2	19315414	34628921	63.77%	4.669%
75	21.327	3114	3118	3123	rVB	20595783	25725653	47.38%	3.469%
76	21.551	3152	3156	3161	rVB	120512	160955	0.30%	0.022%
77	22.857	3370	3378	3382	rBV	14368113	27186069	50.07%	3.666%
78	22.904	3382	3386	3398	rVB	12561080	20002575	36.84%	2.697%
79	23.380	3459	3467	3471	rBV	9129323	18758755	34.55%	2.529%
80	23.433	3471	3476	3483	rVV	12124576	22779233	41.95%	3.072%
81	23.521	3484	3491	3500	rVB	4766116	9260443	17.05%	1.249%
82	24.086	3583	3587	3594	rVB2	243555	423184	0.78%	0.057%
83	24.833	3711	3714	3722	rVB	241634	456655	0.84%	0.062%
84	25.780	3858	3875	3894	rVB3	16255065	54300257	100.00%	7.322%

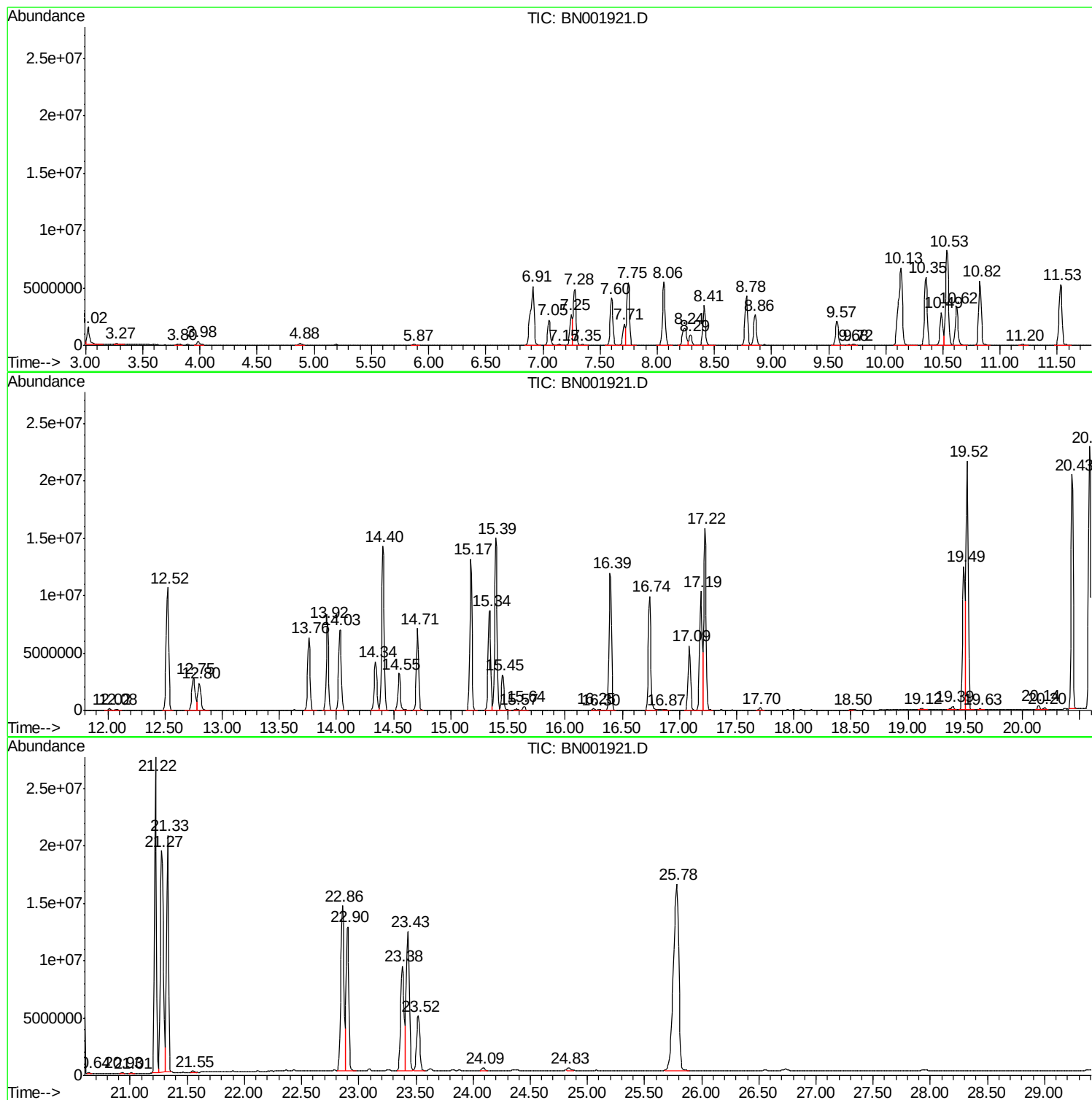
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P



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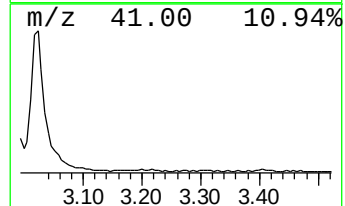
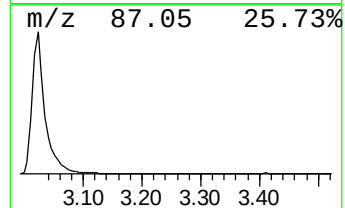
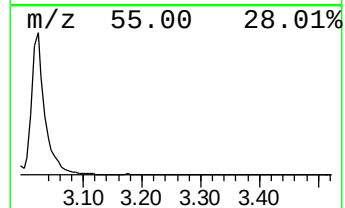
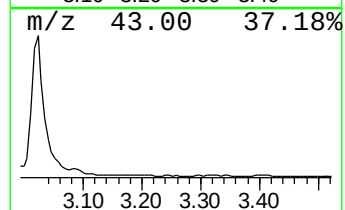
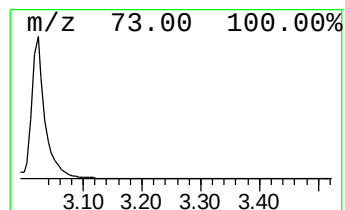
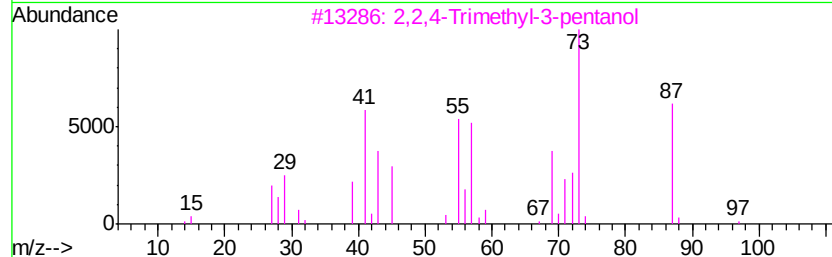
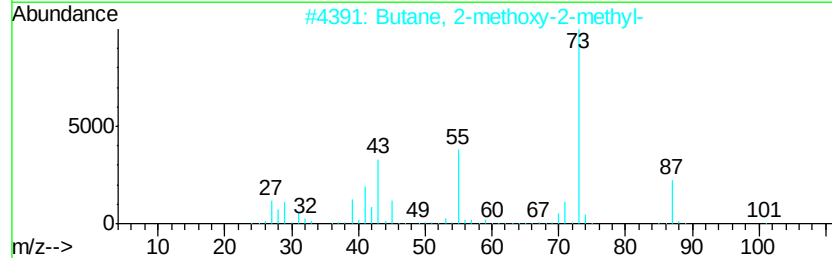
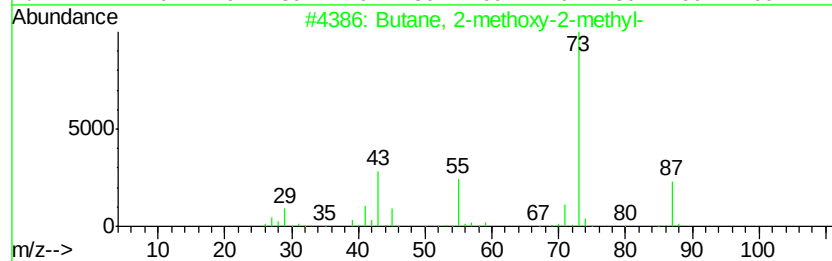
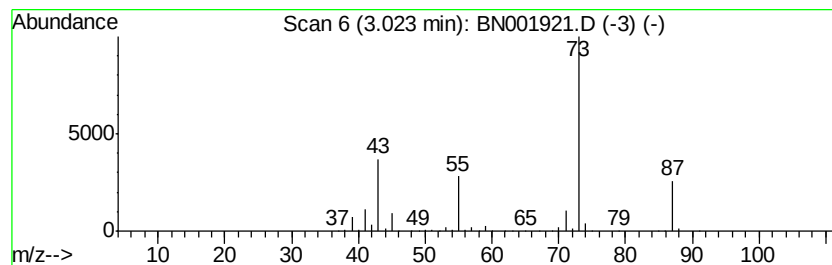
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN061918.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.02	16.33 ng/ul	2425360	1,4-Dichlorobenzene-d4	7.71

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	83
2		Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	78
3		2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	40
4		1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	23
5		Pentane, 3-methoxy-	102	C6H14O	036839-67-5	12



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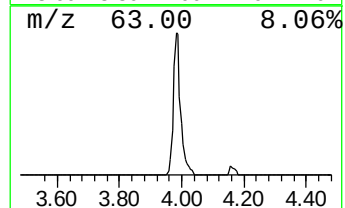
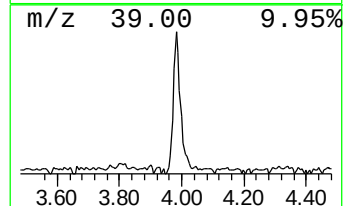
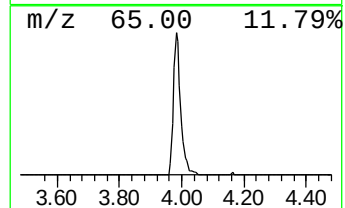
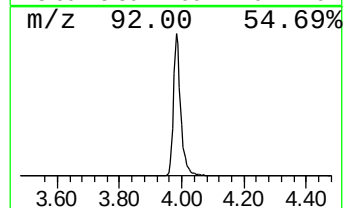
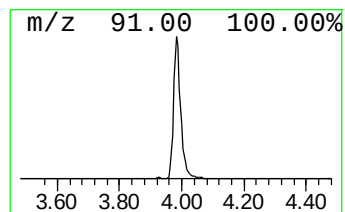
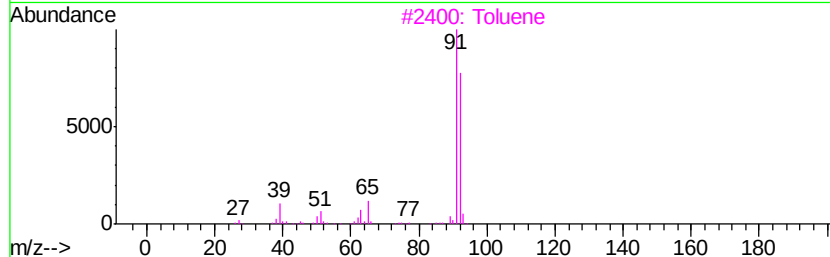
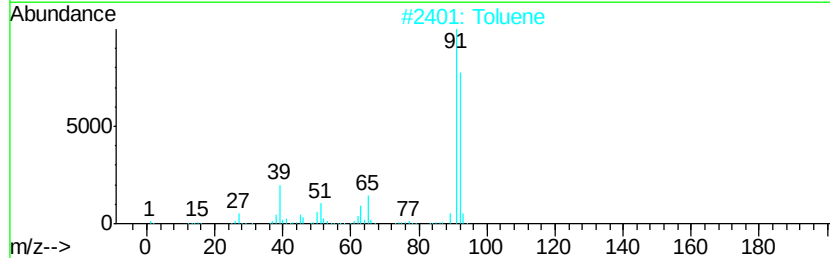
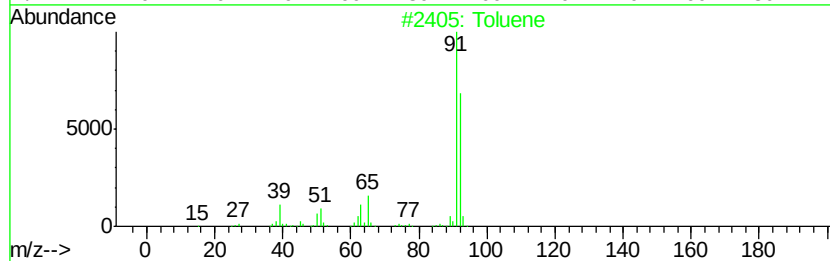
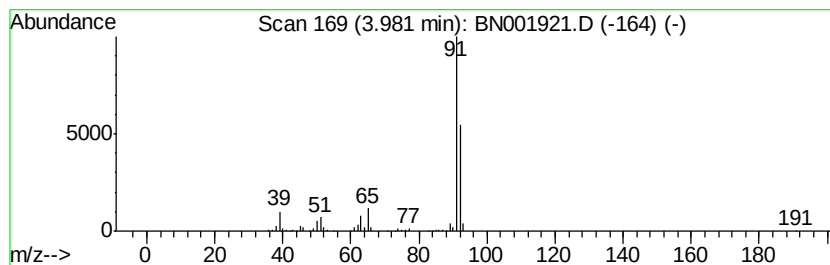
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 Toluene Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.98	3.30 ng/ul	490662	1,4-Dichlorobenzene-d4	7.71

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Toluene			92	C7H8	000108-88-3	95
2	Toluene			92	C7H8	000108-88-3	93
3	Toluene			92	C7H8	000108-88-3	91
4	Toluene			92	C7H8	000108-88-3	91
5	1,3,5-Cycloheptatriene			92	C7H8	000544-25-2	90



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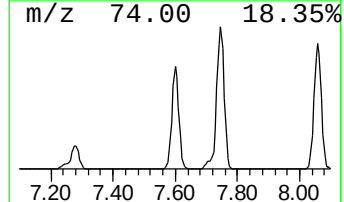
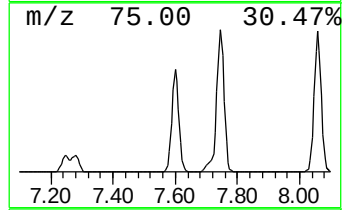
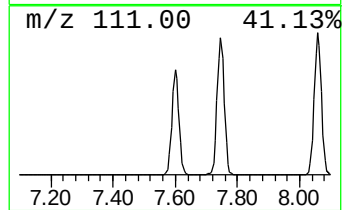
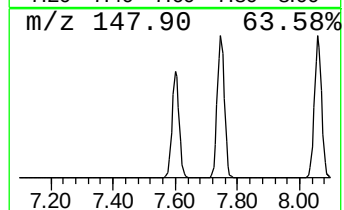
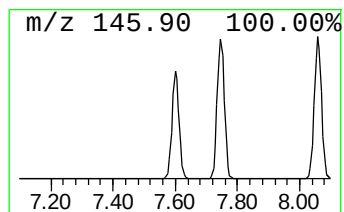
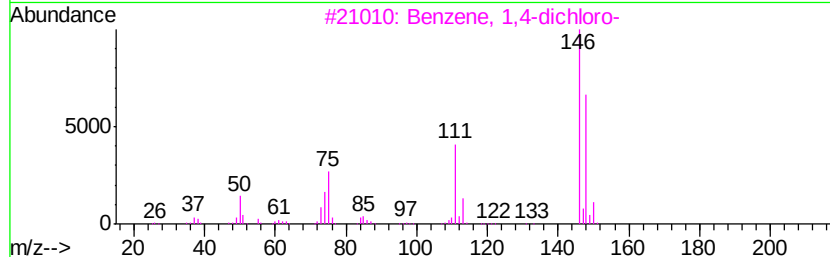
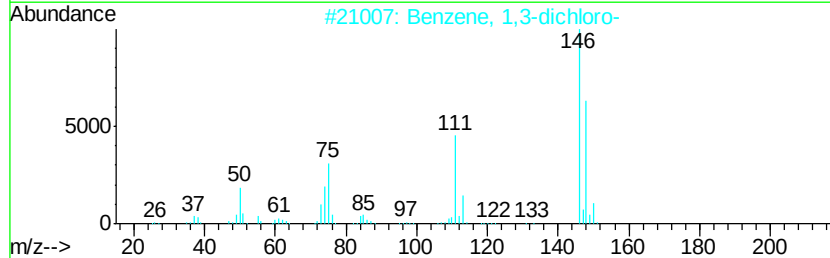
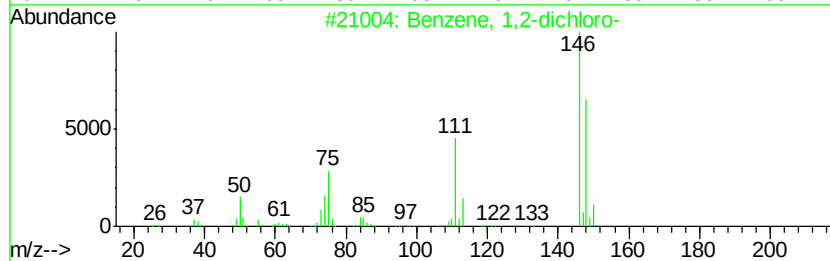
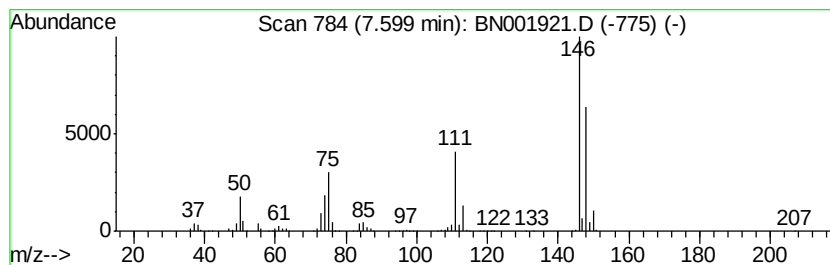
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.60	44.33 ng/ul	6584210	1,4-Dichlorobenzene-d4	7.71

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,2-dichloro-	146	C6H4Cl2	000095-50-1	98
2		Benzene, 1,3-dichloro-	146	C6H4Cl2	000541-73-1	98
3		Benzene, 1,4-dichloro-	146	C6H4Cl2	000106-46-7	98
4		Benzene, 1,3-dichloro-	146	C6H4Cl2	000541-73-1	97
5		Benzene, 1,4-dichloro-	146	C6H4Cl2	000106-46-7	96



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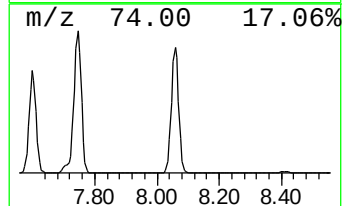
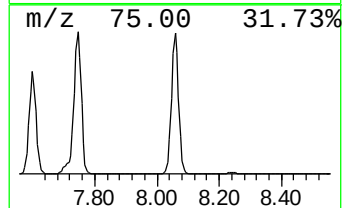
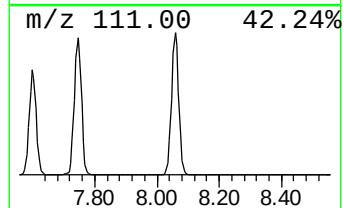
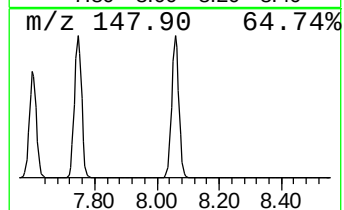
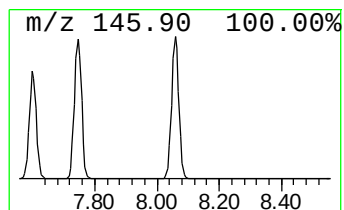
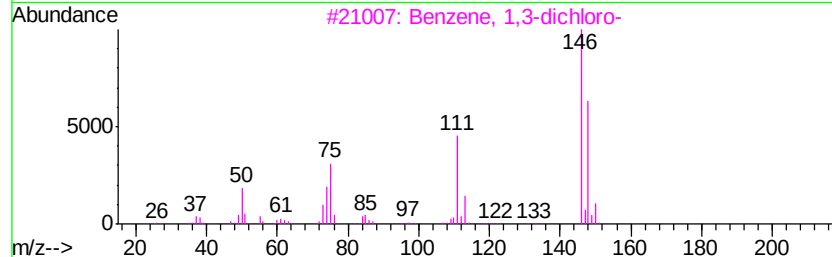
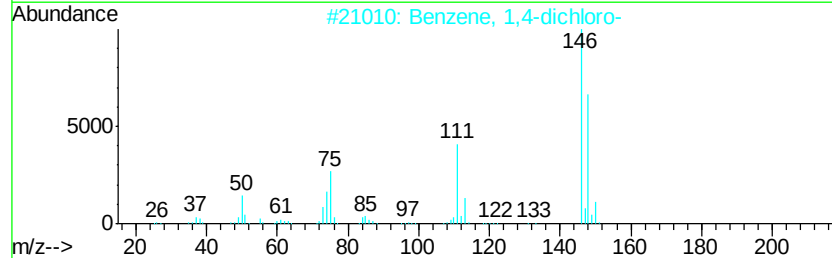
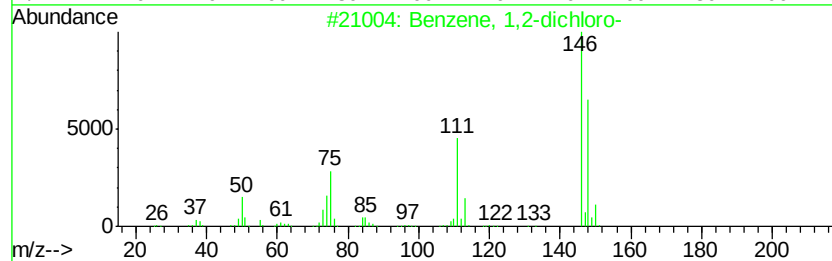
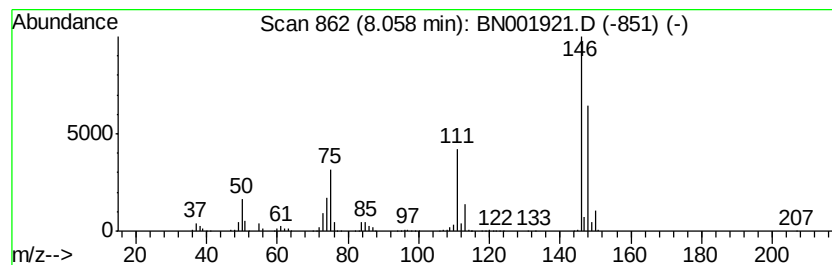
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 Benzene, 1,4-dichloro- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.06	61.01 ng/ul	9061430	1,4-Dichlorobenzene-d4	7.71

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,2-dichloro-	146	C6H4Cl2	000095-50-1	98
2		Benzene, 1,4-dichloro-	146	C6H4Cl2	000106-46-7	98
3		Benzene, 1,3-dichloro-	146	C6H4Cl2	000541-73-1	98
4		Benzene, 1,3-dichloro-	146	C6H4Cl2	000541-73-1	97
5		Benzene, 1,4-dichloro-	146	C6H4Cl2	000106-46-7	97



Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN062918\
Data File : BN001921.D
Acq On : 29 Jun 2018 09:58
Operator : JU/SJ
Sample : J3661-02
Misc :
ALS Vial : 3 Sample Multiplier: 1

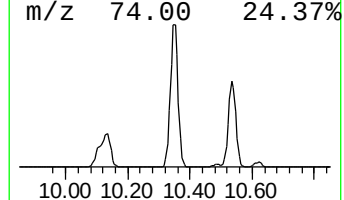
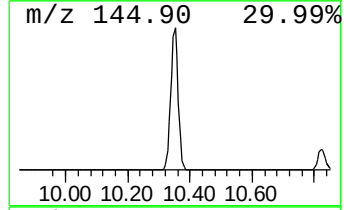
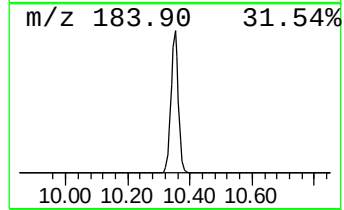
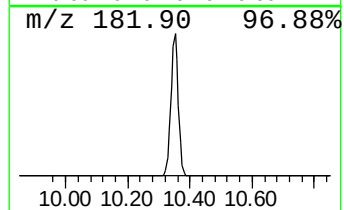
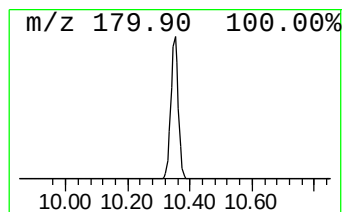
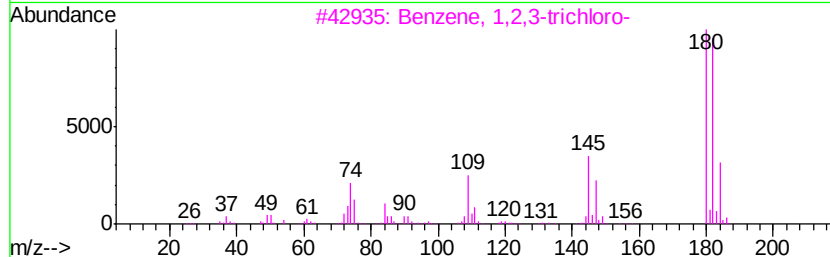
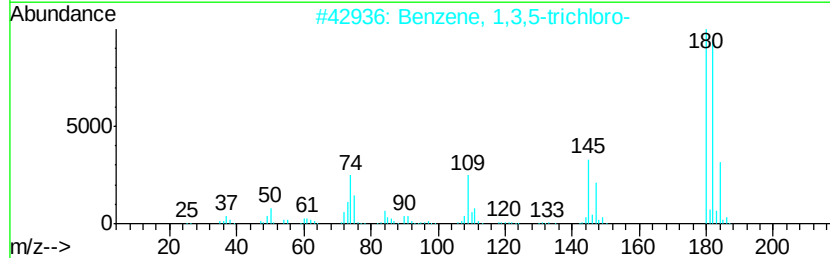
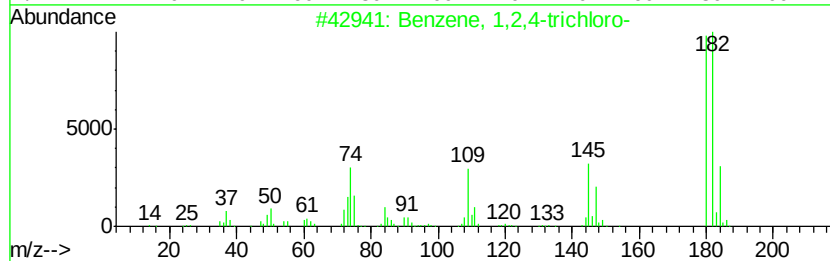
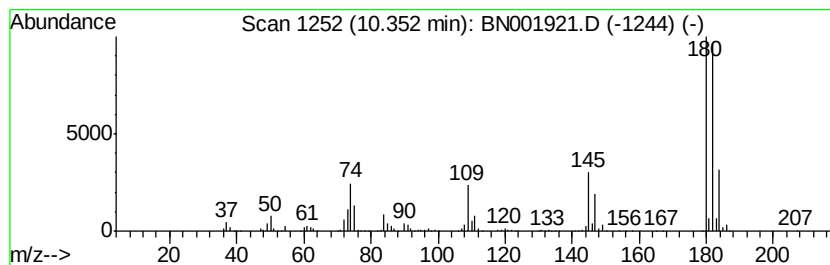
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ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN061918.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 Benzene, 1,2,4-trichloro- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.35	41.47 ng/ul	10007700	Napthalene-d8	10.49
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Benzene, 1,2,4-trichloro-	180 C6H3Cl3	000120-82-1 98
2		Benzene, 1,3,5-trichloro-	180 C6H3Cl3	000108-70-3 98
3		Benzene, 1,2,3-trichloro-	180 C6H3Cl3	000087-61-6 98
4		Benzene, 1,3,5-trichloro-	180 C6H3Cl3	000108-70-3 98
5		Benzene, 1,2,3-trichloro-	180 C6H3Cl3	000087-61-6 98



Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN062918\
Data File : BN001921.D
Acq On : 29 Jun 2018 09:58
Operator : JU/SJ
Sample : J3661-02
Misc :
ALS Vial : 3 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
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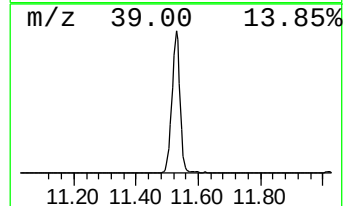
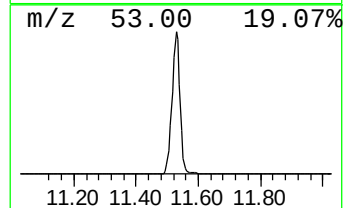
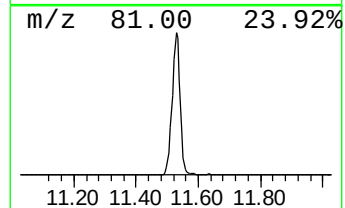
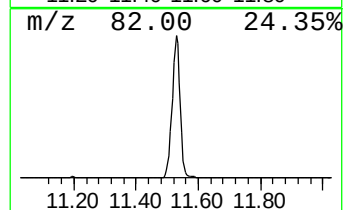
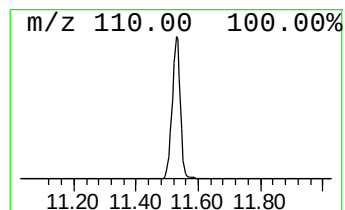
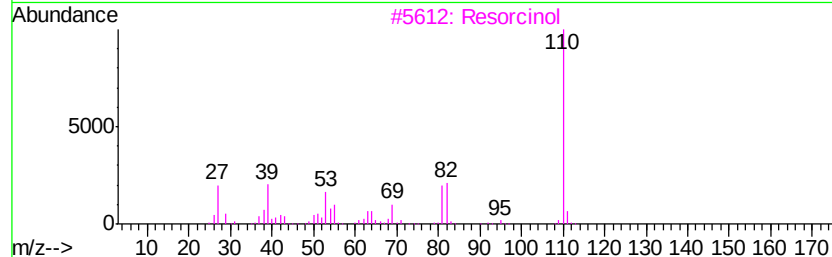
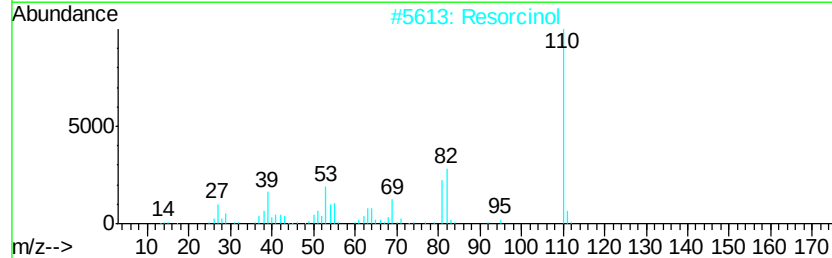
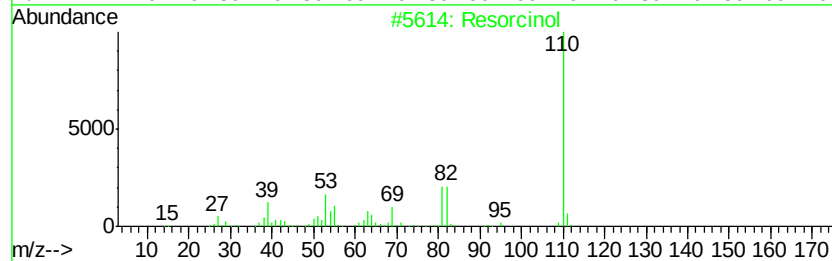
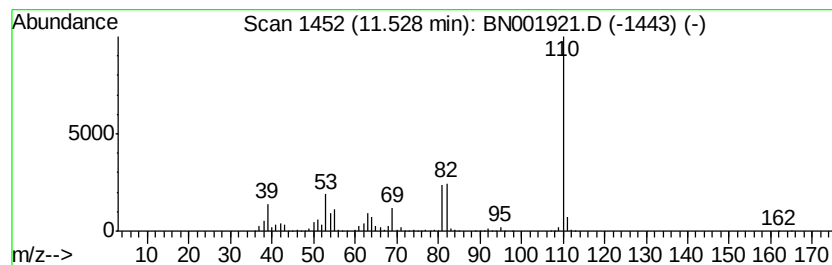
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 7 Resorcinol Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.53	39.75 ng/ul	9592110	Napthalene-d8	10.49

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Resorcinol	110	C6H6O2	000108-46-3	97
2		Resorcinol	110	C6H6O2	000108-46-3	94
3		Resorcinol	110	C6H6O2	000108-46-3	94
4		Resorcinol	110	C6H6O2	000108-46-3	90
5		1,3-Benzenediol, 4,4'-thiobis-	250	C12H10O4S	000097-29-0	83



Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN062918\
Data File : BN001921.D
Acq On : 29 Jun 2018 09:58
Operator : JU/SJ
Sample : J3661-02
Misc :
ALS Vial : 3 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
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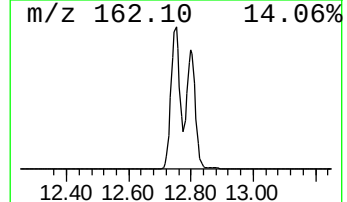
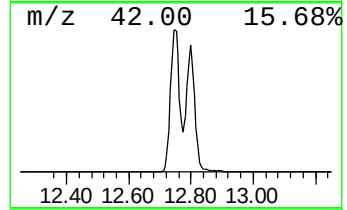
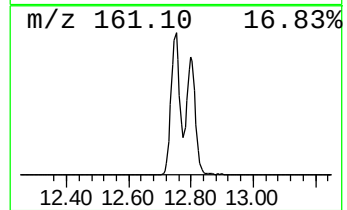
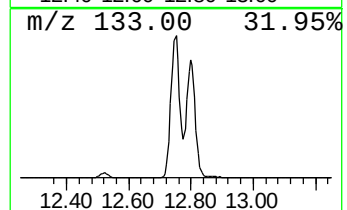
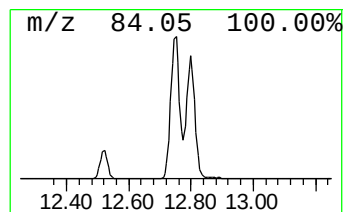
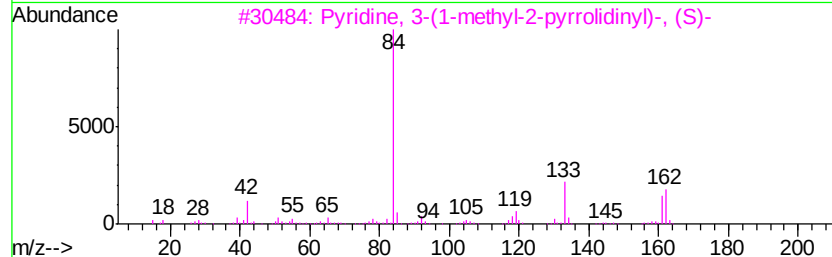
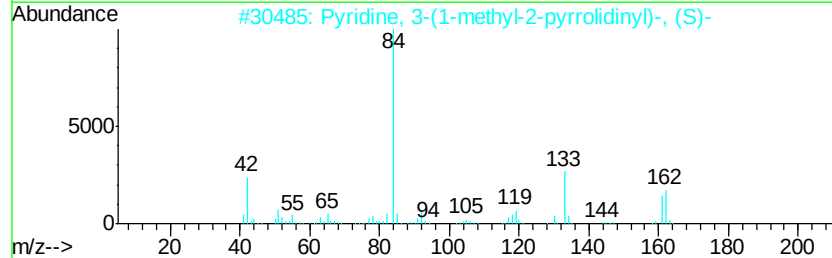
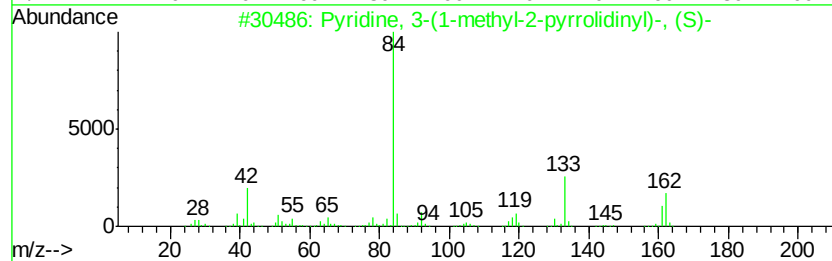
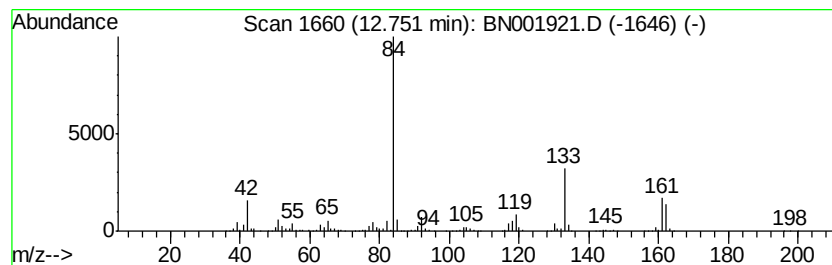
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 8 Pyridine, 3-(1-methyl-2-pyr... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.75	16.91 ng/ul	5653840	Acenaphthene-d10	14.34

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Pyridine, 3-(1-methyl-2-pyrrolid...	162	C10H14N2	000054-11-5	95
2		Pyridine, 3-(1-methyl-2-pyrrolid...	162	C10H14N2	000054-11-5	95
3		Pyridine, 3-(1-methyl-2-pyrrolid...	162	C10H14N2	000054-11-5	90
4		Pyridine, 2-(1-methyl-2-pyrrolid...	162	C10H14N2	023950-04-1	70
5		Pyridine, 3-(1-methyl-2-pyrrolid...	162	C10H14N2	000054-11-5	47



Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN062918\
Data File : BN001921.D
Acq On : 29 Jun 2018 09:58
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Sample : J3661-02
Misc :
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Instrument :
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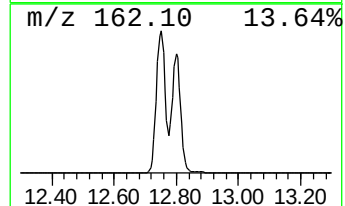
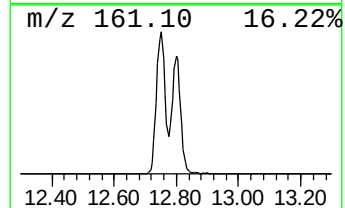
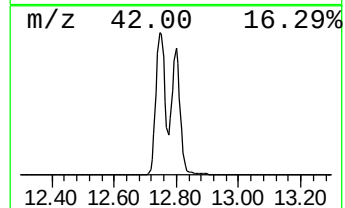
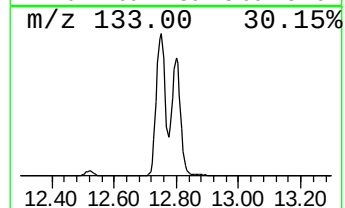
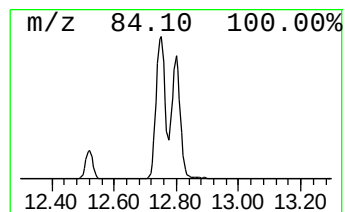
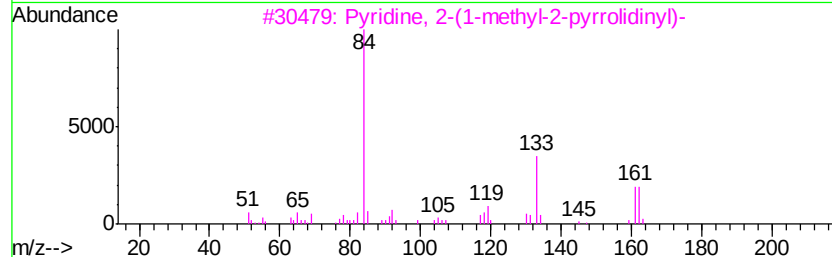
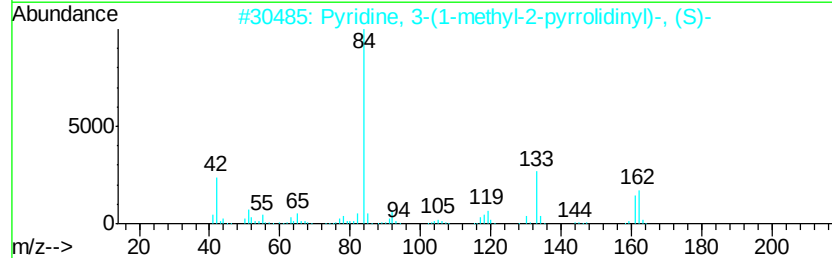
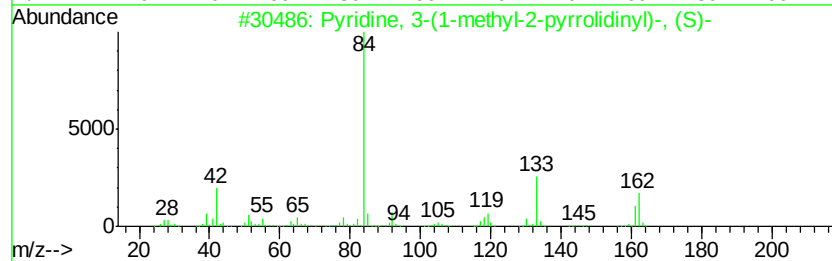
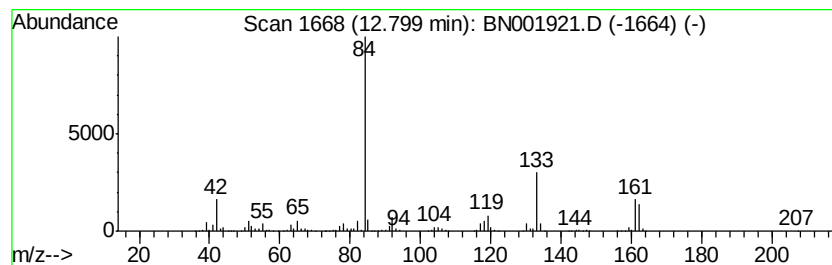
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 9 unknown-01 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.80	13.22 ng/ul	4419720	Acenaphthene-d10	14.34

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Pyridine, 3-(1-methyl-2-pyrrolid...	162	C10H14N2	000054-11-5	96
2		Pyridine, 3-(1-methyl-2-pyrrolid...	162	C10H14N2	000054-11-5	95
3		Pyridine, 2-(1-methyl-2-pyrrolid...	162	C10H14N2	023950-04-1	94
4		Pyridine, 3-(1-methyl-2-pyrrolid...	162	C10H14N2	000054-11-5	87
5		Pyridine, 3-(1-methyl-2-pyrrolid...	162	C10H14N2	000054-11-5	50



Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN062918\
Data File : BN001921.D
Acq On : 29 Jun 2018 09:58
Operator : JU/SJ
Sample : J3661-02
Misc :
ALS Vial : 3 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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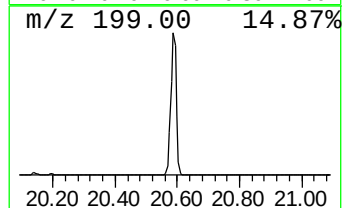
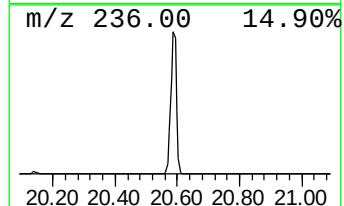
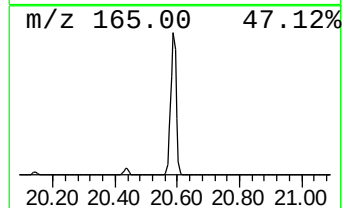
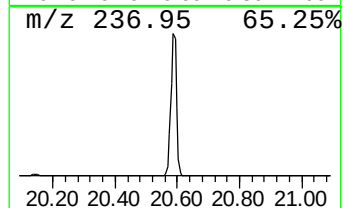
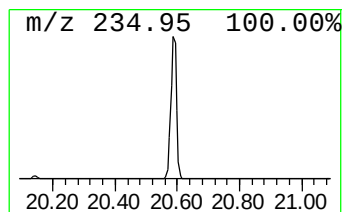
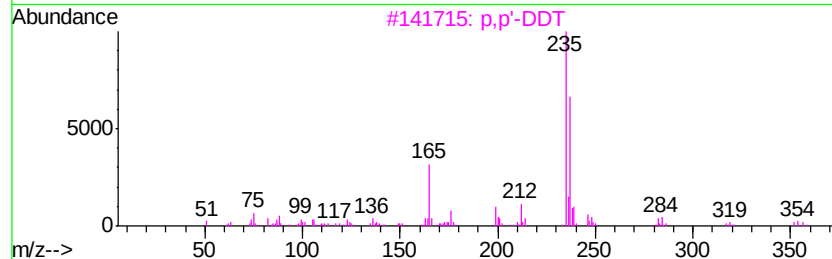
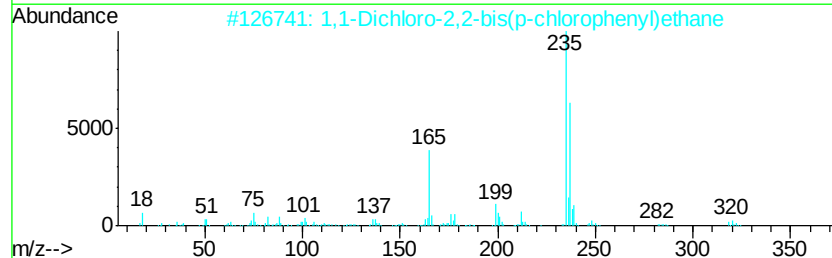
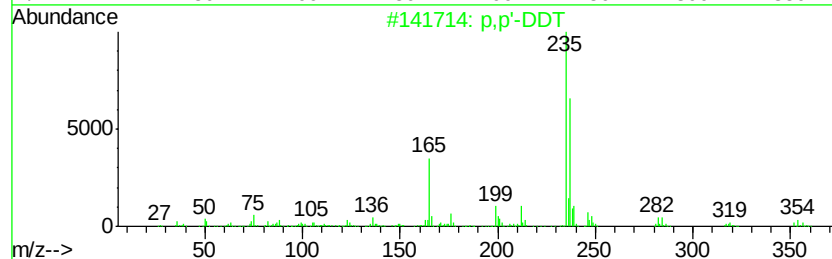
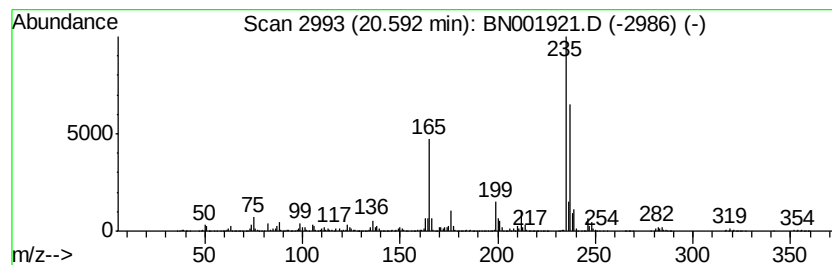
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 10 p,p'-DDT Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.59	16.45 ng/ul	28490300	Chrysene-d12	21.29

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	p,p'-DDT	352	C14H9Cl5	000050-29-3	99
2		1,1-Dichloro-2,2-bis(p-chlorophe...	318	C14H10Cl4	000072-54-8	97
3		p,p'-DDT	352	C14H9Cl5	000050-29-3	95
4		3-Butanone, 1,1-bis(4-chlorophen...	320	C18H18Cl2O	1000158-20-4	93
5		m,p'-DDD	318	C14H10Cl4	004329-12-8	93



Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN062918\
Data File : BN001921.D
Acq On : 29 Jun 2018 09:58
Operator : JU/SJ
Sample : J3661-02
Misc :
ALS Vial : 3 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Butane, 2-methoxy...	3.02	16.3	ng/ul	2425360	1	7.71	2970390	20.0
Toluene	3.98	3.3	ng/ul	490662	1	7.71	2970390	20.0
Benzene, 1,2-dich...	7.60	44.3	ng/ul	6584210	1	7.71	2970390	20.0
Benzene, 1,4-dich...	8.06	61.0	ng/ul	9061430	1	7.71	2970390	20.0
Benzene, 1,2,4-tr...	10.35	41.5	ng/ul	10007700	2	10.49	4826760	20.0
Resorcinol	11.53	39.8	ng/ul	9592110	2	10.49	4826760	20.0
Pyridine, 3-(1-me...	12.75	16.9	ng/ul	5653840	3	14.34	6687800	20.0
unknown-01	12.80	13.2	ng/ul	4419720	3	14.34	6687800	20.0
p,p'-DDT	20.59	16.4	ng/ul	28490300	5	21.29	34628900	20.0