

Data Path : Z:\svoasrv\HPCHEM1\BNA N\Data\BN022420\
 Data File : BN009898.D
 Acq On : 25 Feb 2020 13:17
 Operator : CG/JU
 Sample : L1582-07DL2 30X
 Misc :
 ALS Vial : 36 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 DBDJ9DL2

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-BN021220MA.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	4.969	332	337	343	rVB	14045	20267	0.16%	0.074%
2	5.093	354	358	368	rVB	20070	40118	0.33%	0.146%
3	5.452	413	419	425	rBV3	14117	22584	0.18%	0.082%
4	6.505	593	598	601	rBV3	10296	13001	0.11%	0.047%
5	6.634	615	620	626	rBV2	14789	23436	0.19%	0.085%
6	7.046	684	690	695	rBV	44557	67837	0.55%	0.247%
7	7.387	742	748	760	rBV	376992	605749	4.92%	2.202%
8	7.522	763	771	778	rVB3	28357	62738	0.51%	0.228%
9	7.746	802	809	817	rBV	168587	261424	2.12%	0.950%
10	7.910	830	837	846	rBV	430566	719298	5.84%	2.614%
11	8.552	941	946	951	rBV6	16400	29474	0.24%	0.107%
12	8.604	951	955	962	rVB2	14439	22328	0.18%	0.081%
13	8.722	969	975	984	rBV	20789	35781	0.29%	0.130%
14	8.846	989	996	1003	rVV3	43971	103950	0.84%	0.378%
15	8.904	1003	1006	1015	rVV3	14882	29701	0.24%	0.108%
16	8.999	1015	1022	1027	rVV5	6033	12541	0.10%	0.046%
17	9.281	1061	1070	1073	rBV7	13954	26438	0.21%	0.096%
18	9.399	1086	1090	1095	rBV	14643	23278	0.19%	0.085%
19	9.552	1106	1116	1127	rVV7	62786	151803	1.23%	0.552%
20	9.646	1127	1132	1134	rVV5	14210	20586	0.17%	0.075%
21	9.681	1134	1138	1147	rVB2	51996	85920	0.70%	0.312%
22	10.134	1208	1215	1218	rBV	509620	871365	7.08%	3.167%
23	10.187	1218	1224	1234	rVV	7632919	12309997	100.00%	44.742%
24	10.299	1237	1243	1256	rVB	221386	379838	3.09%	1.381%
25	11.693	1475	1480	1486	rBV3	18920	29746	0.24%	0.108%
26	11.804	1491	1499	1513	rBV	926739	1524969	12.39%	5.543%
27	11.928	1515	1520	1527	rVV	20950	42253	0.34%	0.154%
28	12.028	1527	1537	1552	rVB	465474	800344	6.50%	2.909%
29	12.851	1671	1677	1691	rBV2	135764	208394	1.69%	0.757%
30	13.051	1706	1711	1719	rBV	27023	45105	0.37%	0.164%
31	13.192	1729	1735	1736	rBV2	25158	35216	0.29%	0.128%
32	13.345	1753	1761	1766	rBV2	43438	77028	0.63%	0.280%
33	13.404	1766	1771	1776	rVV3	25009	36439	0.30%	0.132%
34	13.463	1776	1781	1786	rVB	25646	38205	0.31%	0.139%

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35	13.587	1798	1802	1805	rBV3	14805	21003	0.17%	0.076%
36	13.716	1819	1824	1828	rBV	30246	44195	0.36%	0.161%
37	13.751	1828	1830	1835	rVB3	9848	12691	0.10%	0.046%
38	14.028	1870	1877	1882	rBV	706844	1073248	8.72%	3.901%
39	14.092	1882	1888	1898	rVV	642021	923591	7.50%	3.357%
40	14.181	1898	1903	1914	rVB	73446	119723	0.97%	0.435%
41	14.434	1940	1946	1953	rBV	314127	457120	3.71%	1.661%
42	14.492	1953	1956	1962	rVB2	19452	31441	0.26%	0.114%
43	15.034	2041	2048	2053	rBV2	41402	61625	0.50%	0.224%
44	15.092	2053	2058	2064	rVV	216022	308355	2.50%	1.121%
45	15.257	2083	2086	2090	rVB3	10197	12635	0.10%	0.046%
46	15.416	2106	2113	2124	rVB6	20960	40844	0.33%	0.148%
47	15.551	2131	2136	2142	rBV5	10201	21133	0.17%	0.077%
48	15.781	2168	2175	2188	rBV3	111425	236667	1.92%	0.860%
49	15.981	2203	2209	2216	rBV4	14654	23446	0.19%	0.085%
50	16.316	2259	2266	2272	rBV5	18093	35649	0.29%	0.130%
51	16.451	2283	2289	2296	rBV7	12153	28271	0.23%	0.103%
52	16.610	2312	2316	2318	rVB2	12732	15888	0.13%	0.058%
53	16.786	2339	2346	2350	rBV	885754	1294529	10.52%	4.705%
54	16.828	2350	2353	2359	rVB	204382	270511	2.20%	0.983%
55	16.886	2359	2363	2369	rBV2	51673	78728	0.64%	0.286%
56	17.198	2408	2416	2424	rBV	354576	506395	4.11%	1.841%
57	17.375	2441	2446	2454	rBV2	10663	18196	0.15%	0.066%
58	17.716	2500	2504	2510	rBV3	19881	26009	0.21%	0.095%
59	17.863	2525	2529	2534	rVB6	8503	13259	0.11%	0.048%
60	18.110	2568	2571	2580	rVB4	4691	12513	0.10%	0.045%
61	19.198	2750	2756	2764	rBV2	55863	80848	0.66%	0.294%
62	20.751	3017	3020	3023	rBV5	15081	20553	0.17%	0.075%
63	20.804	3026	3029	3036	rVV2	29181	46123	0.37%	0.168%
64	21.004	3058	3063	3073	rBV2	1205650	1466650	11.91%	5.331%
65	23.104	3414	3420	3431	rVB	820189	1434419	11.65%	5.214%

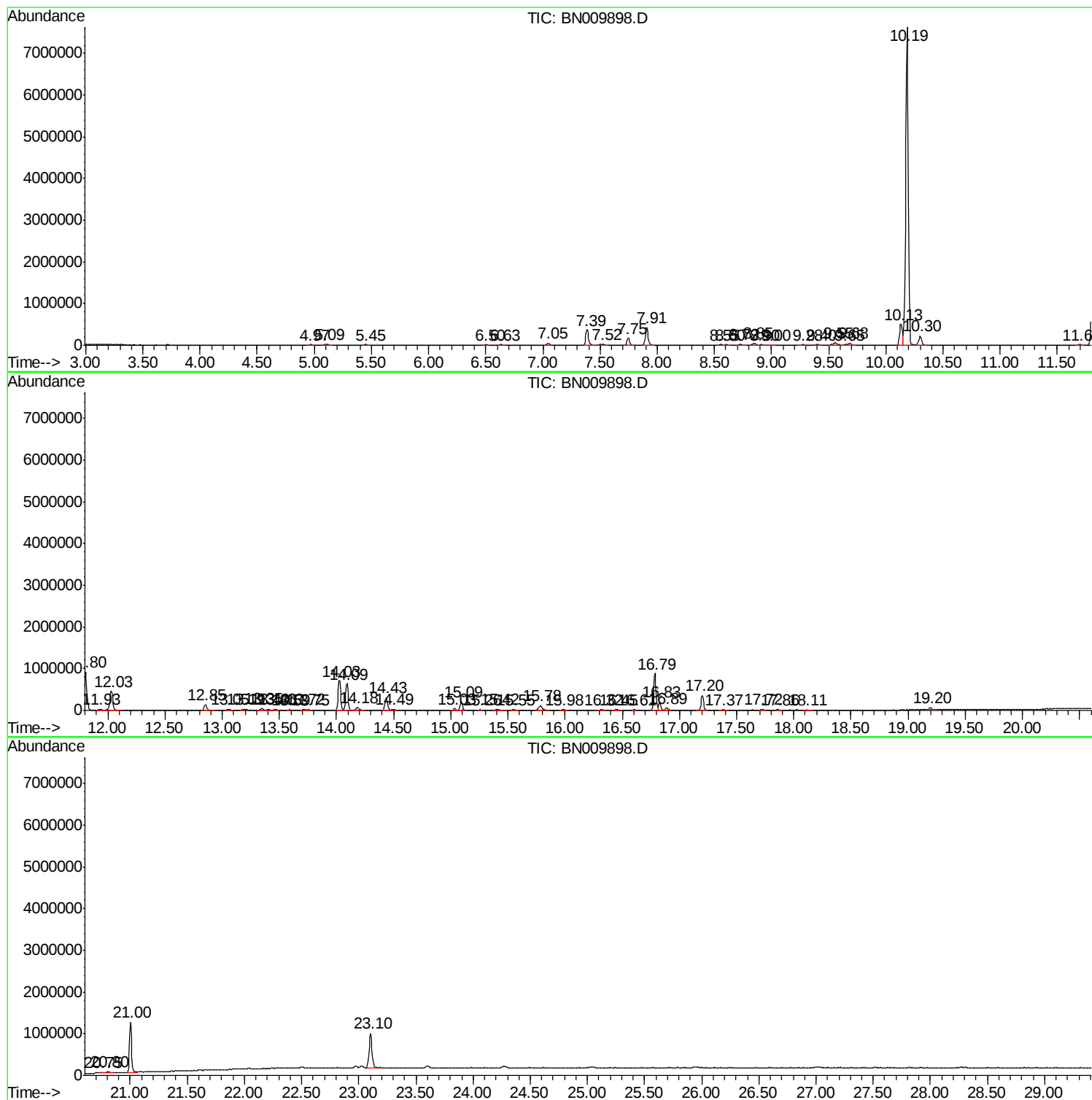
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Quant Title : SVOA CALIBRATION

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



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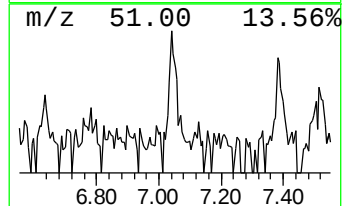
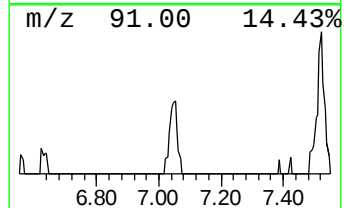
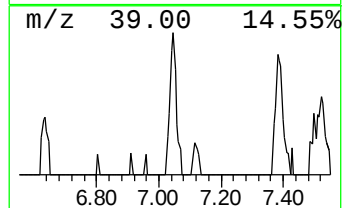
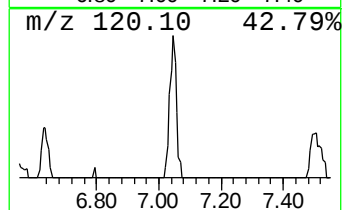
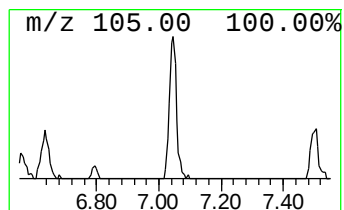
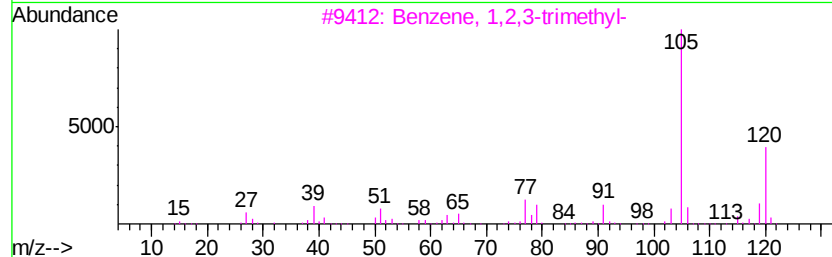
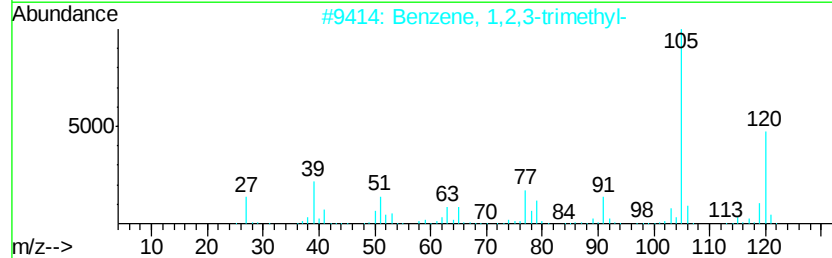
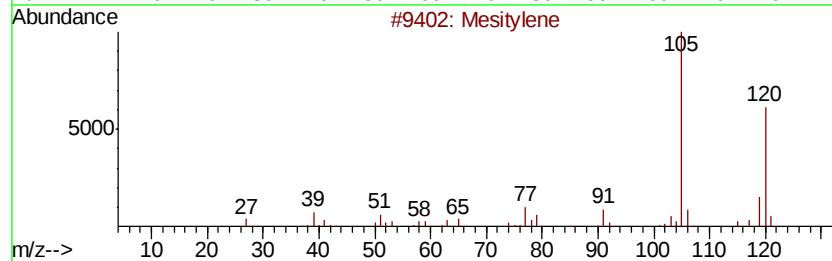
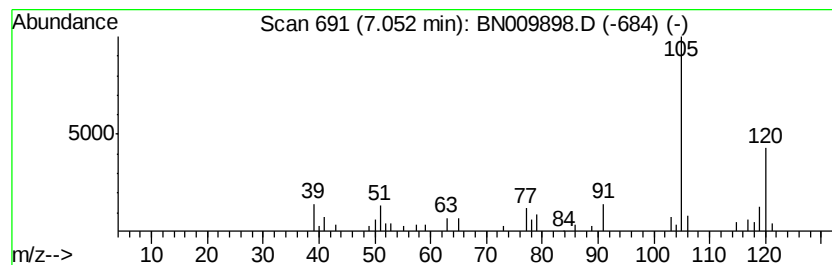
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Quant Title : SVOA CALIBRATION

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

Peak Number 1 Mesitylene Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.05	2.24 ng/ul	67837	1,4-Dichlorobenzene-d4	7.39

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Mesitylene	120	C9H12	000108-67-8	95
2		Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	94
3		Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	94
4		Mesitylene	120	C9H12	000108-67-8	94
5		Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	94



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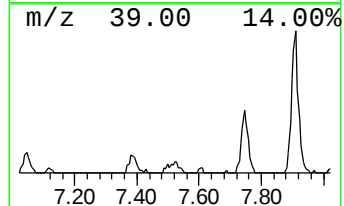
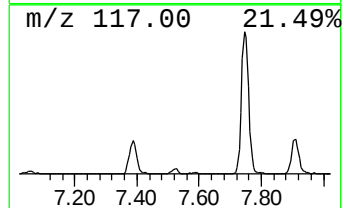
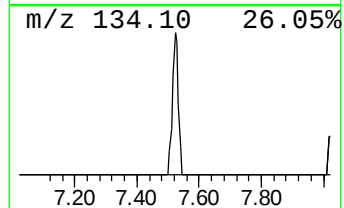
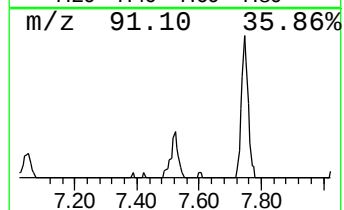
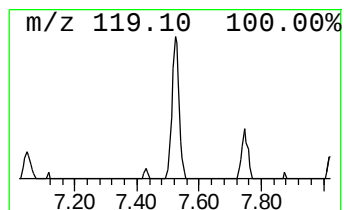
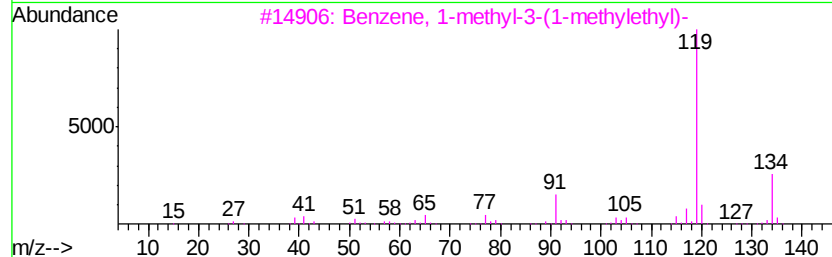
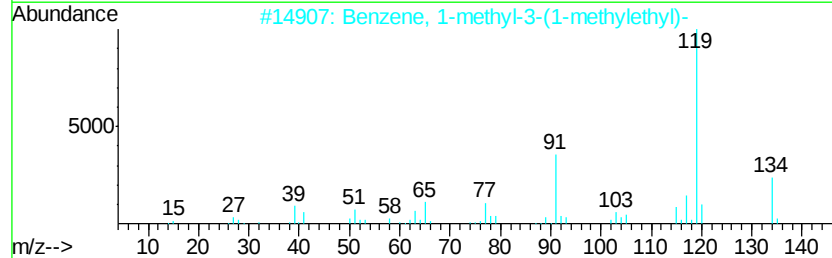
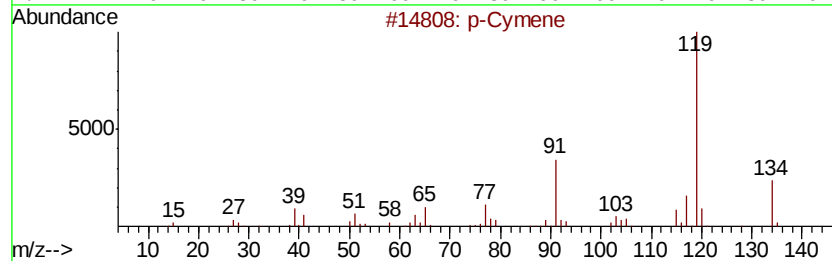
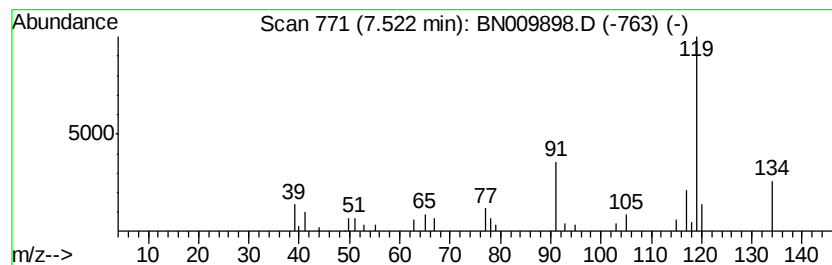
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN021220MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 p-Cymene Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.52	2.07 ng/ul	62738	1,4-Dichlorobenzene-d4	7.39

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	p-Cymene	134	C10H14	000099-87-6	80
2		Benzene, 1-methyl-3-(1-methylethyl)-	134	C10H14	000535-77-3	78
3		Benzene, 1-methyl-3-(1-methylethyl)-	134	C10H14	000535-77-3	72
4		Benzene, 1-methyl-3-(1-methylethyl)-	134	C10H14	000535-77-3	64
5		o-Cymene	134	C10H14	000527-84-4	64



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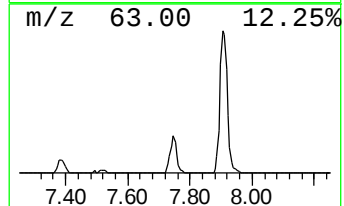
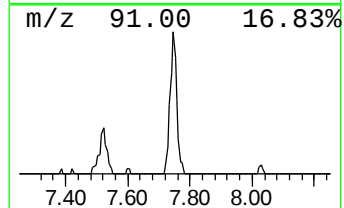
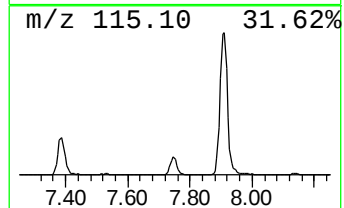
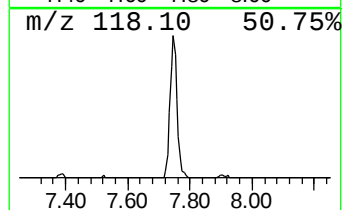
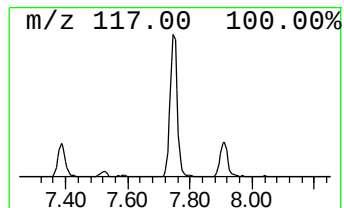
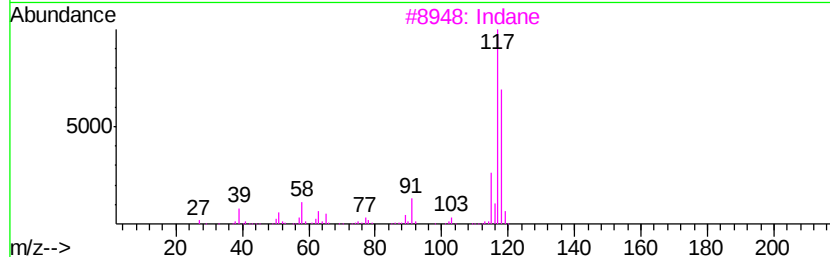
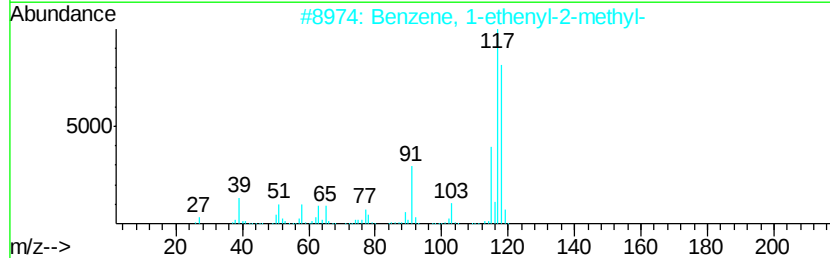
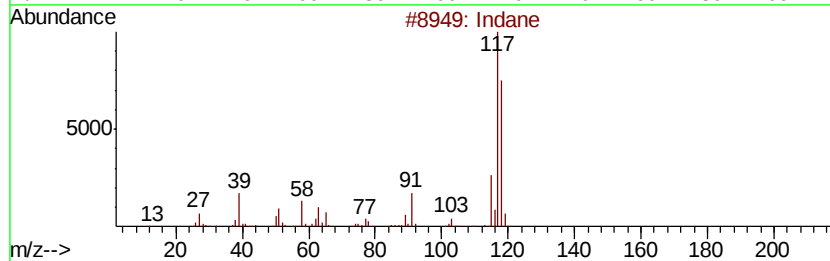
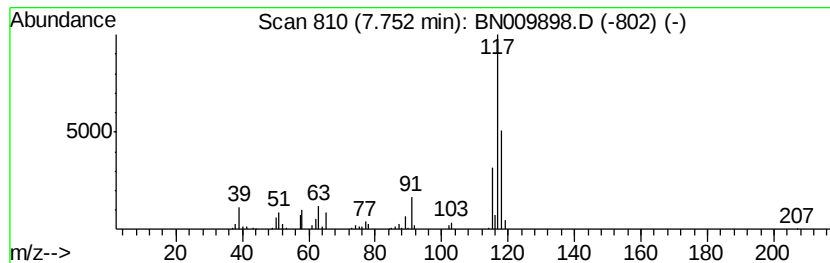
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN021220MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 Indane Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.75	8.63 ng/ul	261424	1,4-Dichlorobenzene-d4	7.39

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Indane	118	C9H10	000496-11-7	95
2		Benzene, 1-ethenyl-2-methyl-	118	C9H10	000611-15-4	81
3		Indane	118	C9H10	000496-11-7	81
4		Benzene, 2-propenyl-	118	C9H10	000300-57-2	76
5		Benzene, 2-propenyl-	118	C9H10	000300-57-2	76



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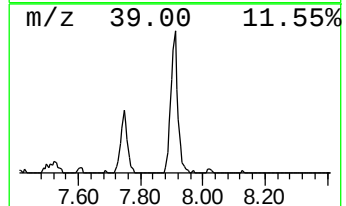
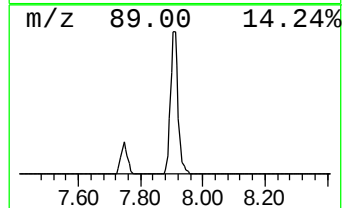
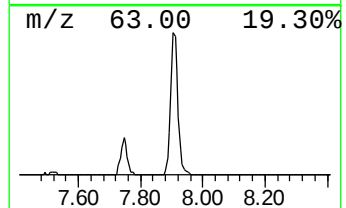
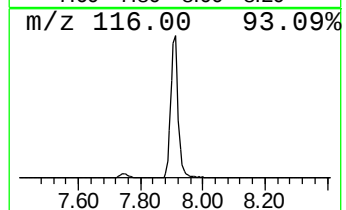
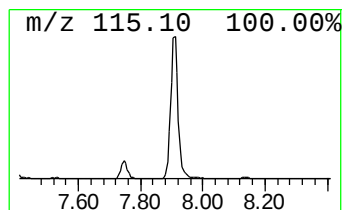
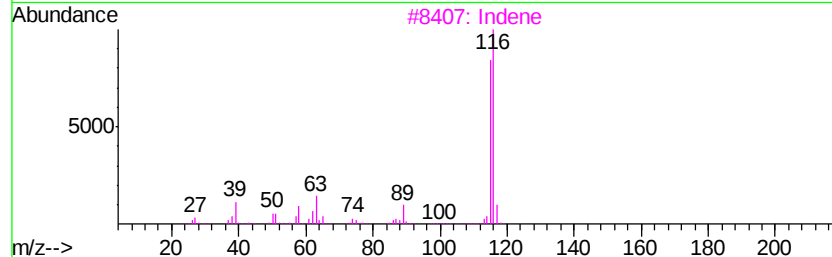
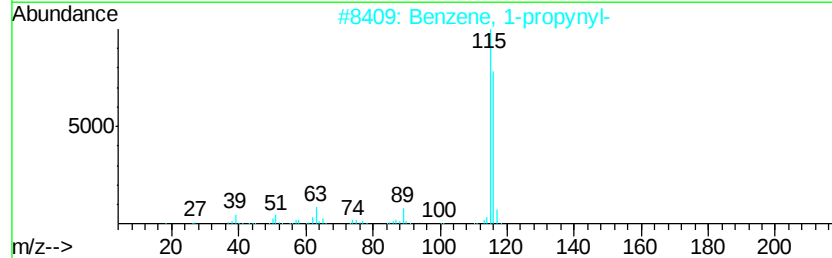
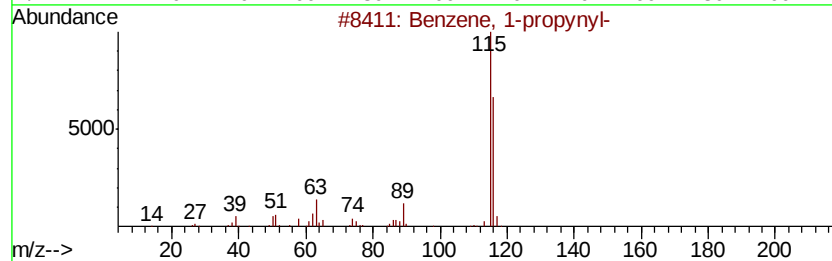
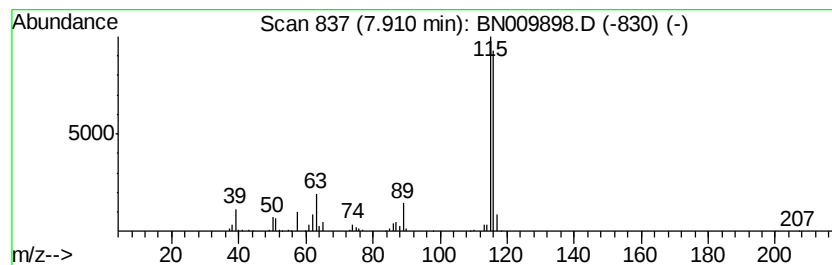
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN021220MA.M
Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.91	23.75 ng/ul	719298	1,4-Dichlorobenzene-d4	7.39

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-propynyl-	116	C9H8	000673-32-5	94
2		Benzene, 1-propynyl-	116	C9H8	000673-32-5	90
3		Indene	116	C9H8	000095-13-6	90
4		Benzene, 1-ethynyl-4-methyl-	116	C9H8	000766-97-2	90
5		Benzene, 1-propynyl-	116	C9H8	000673-32-5	87



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DBDJ9DL2

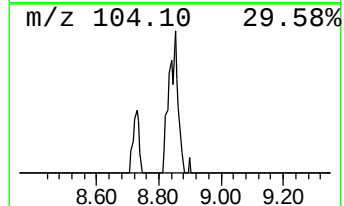
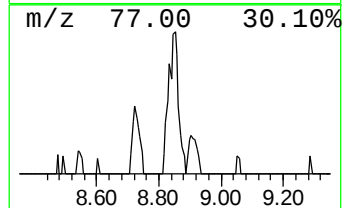
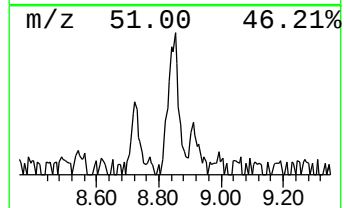
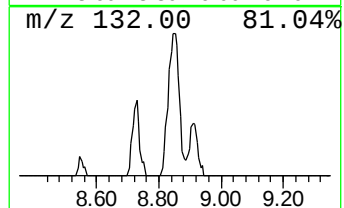
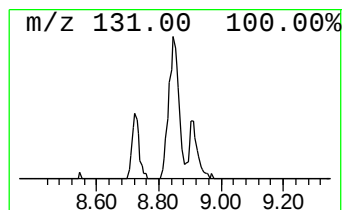
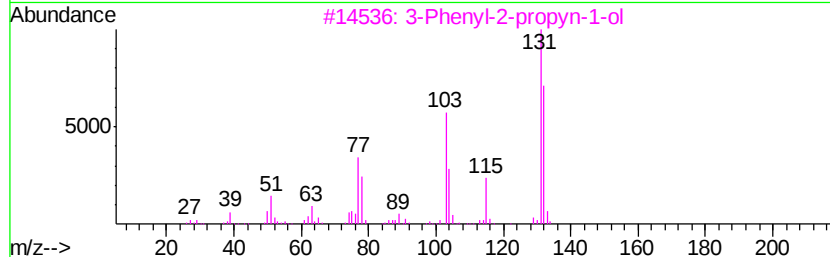
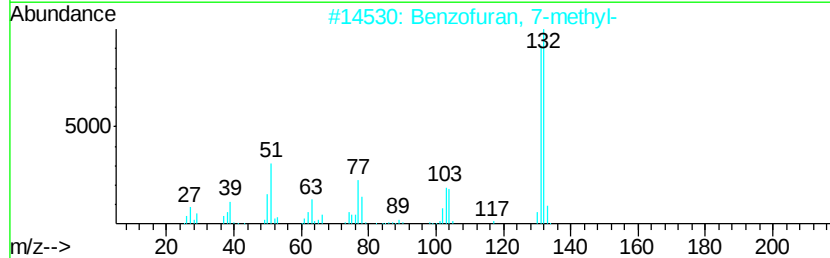
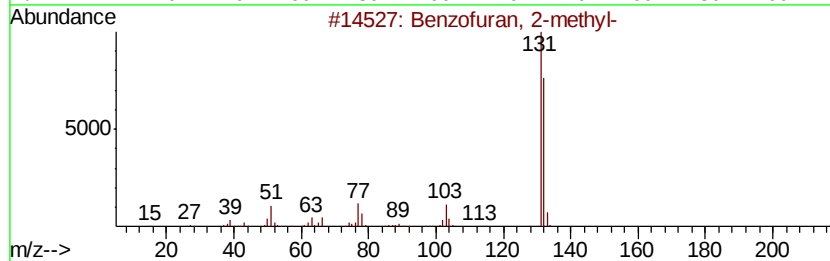
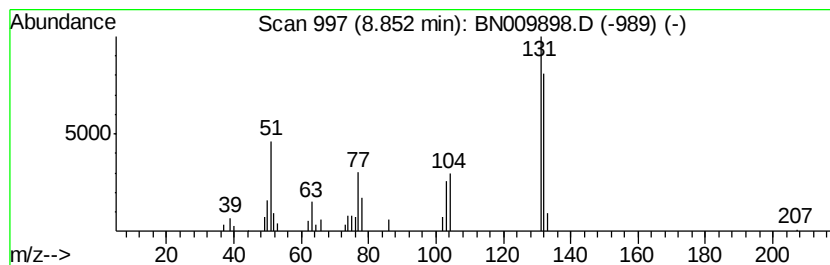
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 Benzofuran, 2-methyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.85	2.39 ng/ul	103950	Naphthalene-d8	10.13

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Benzofuran, 2-methyl-	132	C9H8O	004265-25-2	94
2		Benzofuran, 7-methyl-	132	C9H8O	017059-52-8	93
3		3-Phenyl-2-propyn-1-ol	132	C9H8O	001504-58-1	91
4		Cinnamaldehyde, (E)-	132	C9H8O	014371-10-9	91
5		2-Propenal, 3-phenyl-	132	C9H8O	000104-55-2	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN022420\
Data File : BN009898.D
Acq On : 25 Feb 2020 13:17
Operator : CG/JU
Sample : L1582-07DL2 30X
Misc :
ALS Vial : 36 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
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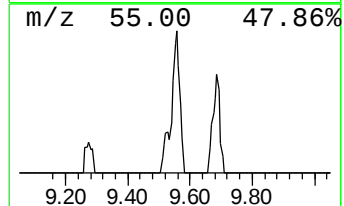
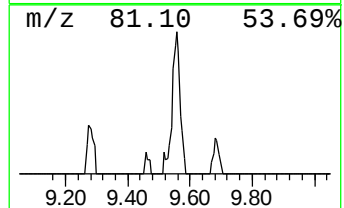
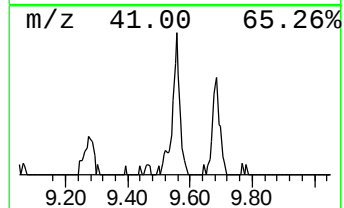
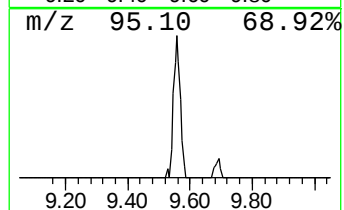
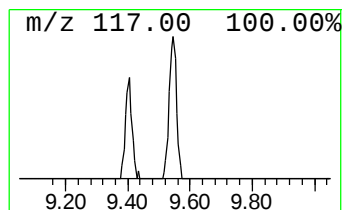
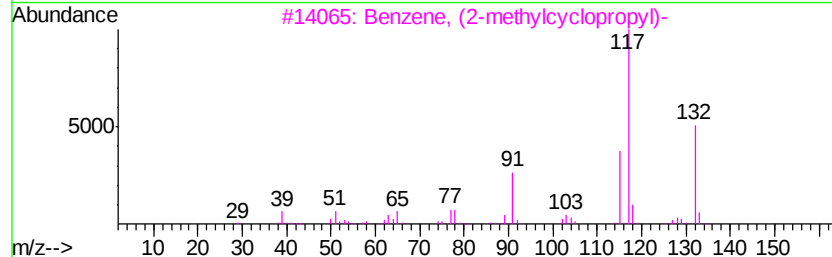
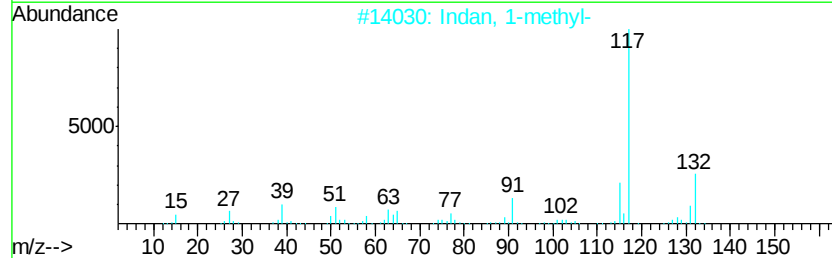
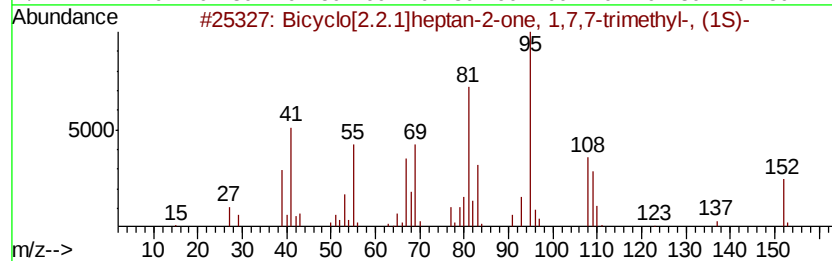
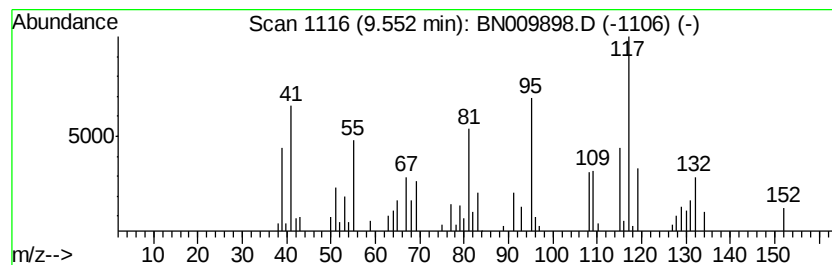
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 Bicyclo[2.2.1]heptan-2-one,... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.55	3.48 ng/ul	151803	Naphthalene-d8	10.13

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Bicyclo[2.2.1]heptan-2-one, 1,7,...	152	C10H16O	000464-48-2	60
2		Indan, 1-methyl-	132	C10H12	000767-58-8	38
3		Benzene, (2-methylcyclopropyl)-	132	C10H12	1000327-39-0	38
4		1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	35
5		Benzene, (2-bromocyclopropyl)-	196	C9H9Br	036617-02-4	35



Data Path : Z:\svoasrv\HPCHEM1\BNA N\Data\BN022420\
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Operator : CG/JU
Sample : L1582-07DL2 30X
Misc :
ALS Vial : 36 Sample Multiplier: 1

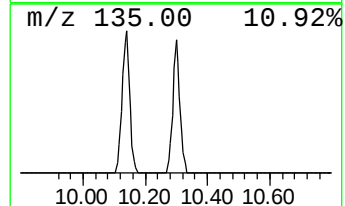
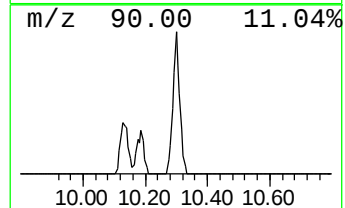
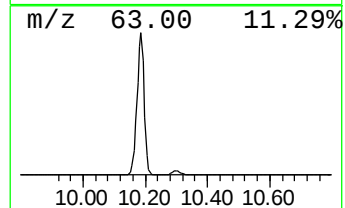
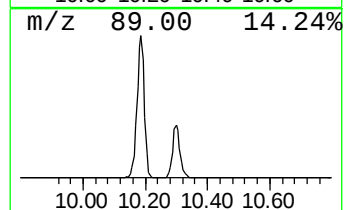
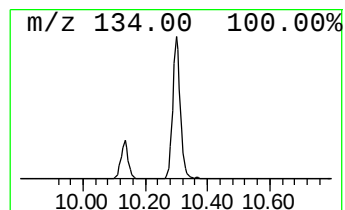
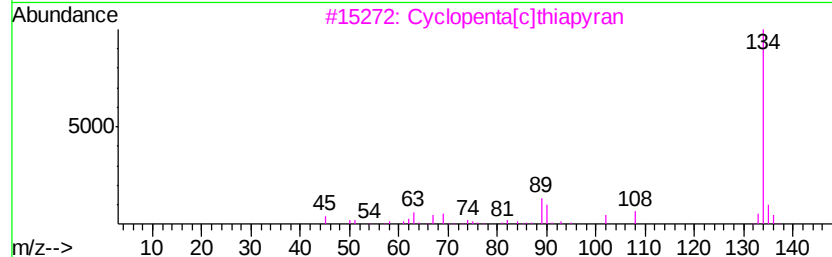
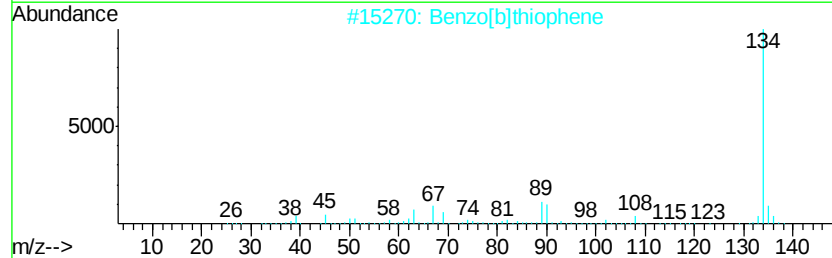
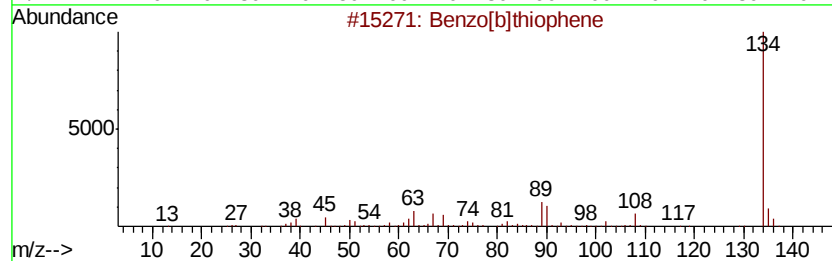
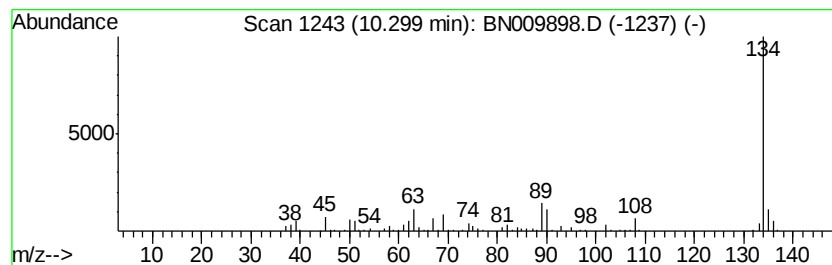
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 Benzo[b]thiophene Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.30	8.72 ng/ul	379838	Napthalene-d8	10.13
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Benzo[b]thiophene	134 C8H6S	000095-15-8	97
2	Benzo[b]thiophene	134 C8H6S	000095-15-8	94
3	Cyclopenta[c]thiapyran	134 C8H6S	000270-63-3	91
4	Benzo[c]thiophene	134 C8H6S	000270-82-6	91
5	Benzo[b]thiophene	134 C8H6S	000095-15-8	90



Data Path : Z:\svoasrv\HPCHEM1\BNA N\Data\BN022420\
Data File : BN009898.D
Acq On : 25 Feb 2020 13:17
Operator : CG/JU
Sample : L1582-07DL2 30X
Misc :
ALS Vial : 36 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
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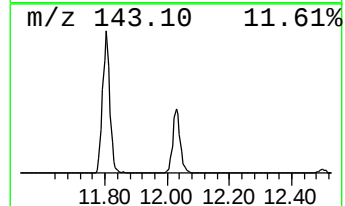
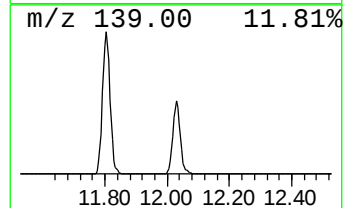
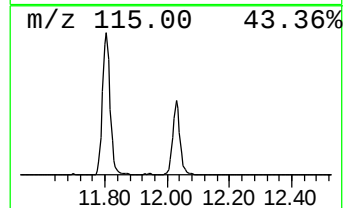
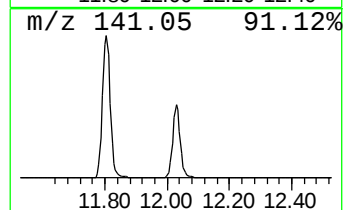
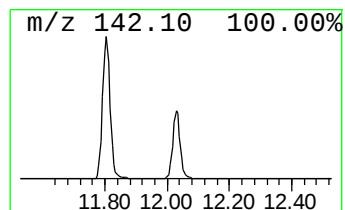
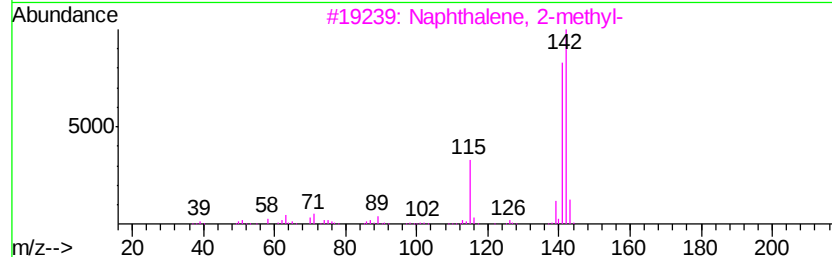
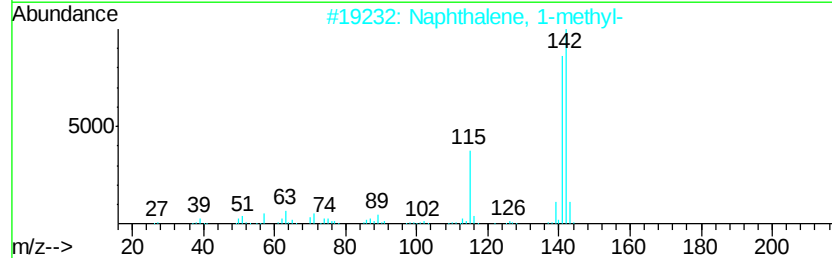
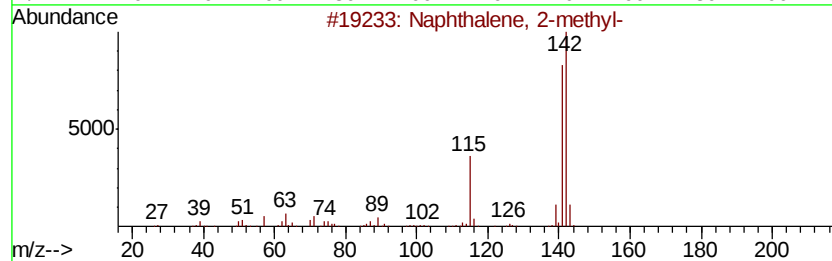
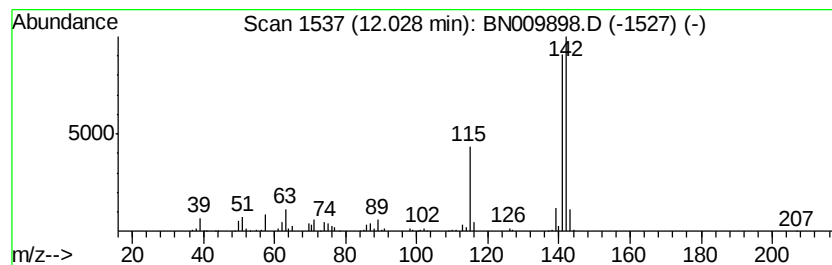
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 Naphthalene, 1-methyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.03	18.37 ng/ul	800344	Naphthalene-d8	10.13

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 2-methyl-	142	C11H10	000091-57-6	96
2		Naphthalene, 1-methyl-	142	C11H10	000090-12-0	96
3		Naphthalene, 2-methyl-	142	C11H10	000091-57-6	95
4		Naphthalene, 1-methyl-	142	C11H10	000090-12-0	94
5		1,4-Methanonaphthalene, 1,4-dihy...	142	C11H10	004453-90-1	94



Data Path : Z:\svoasrv\HPCHEM1\BNA N\Data\BN022420\
Data File : BN009898.D
Acq On : 25 Feb 2020 13:17
Operator : CG/JU
Sample : L1582-07DL2 30X
Misc :
ALS Vial : 36 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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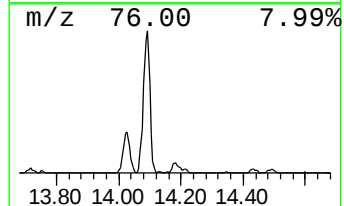
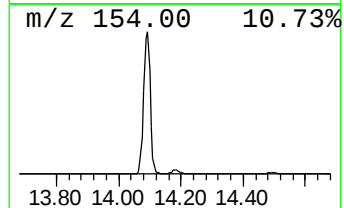
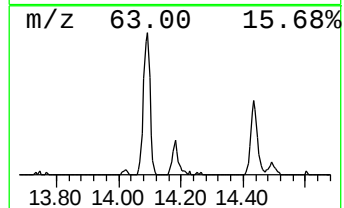
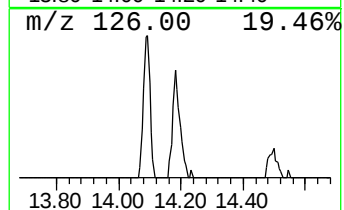
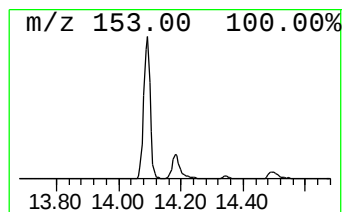
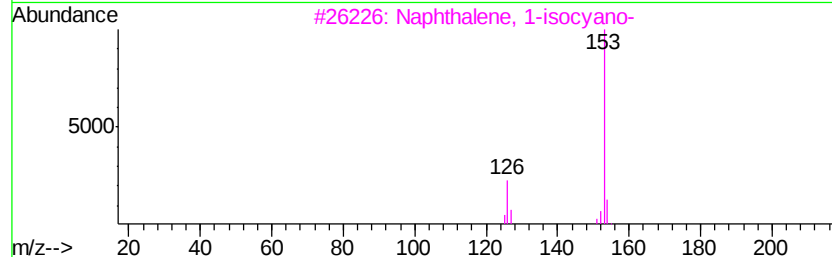
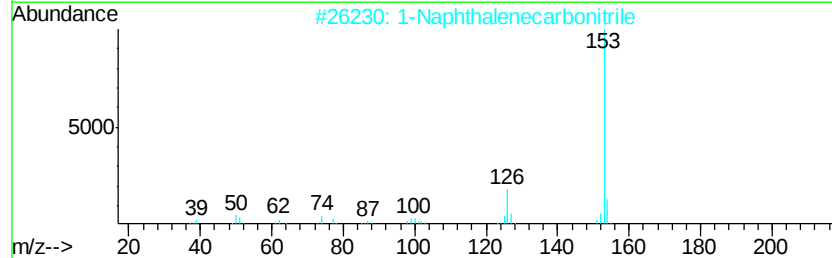
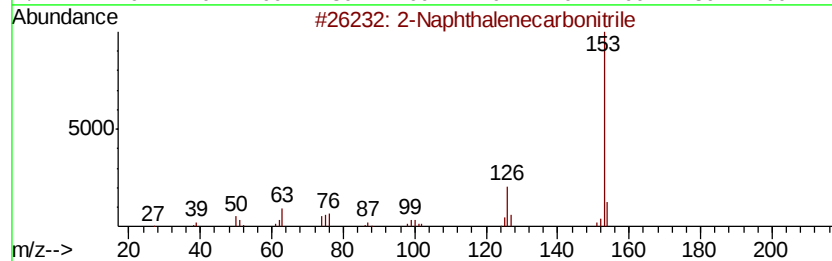
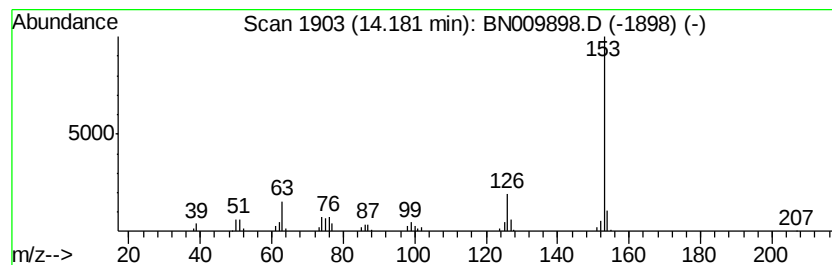
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 2-Naphthalenecarbonitrile Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.18	2.23 ng/ul	119723	Acenaphthene-d10	14.03

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Naphthalenecarbonitrile	153	C11H7N	000613-46-7	97
2		1-Naphthalenecarbonitrile	153	C11H7N	000086-53-3	93
3		Naphthalene, 1-isocvano-	153	C11H7N	001984-04-9	90
4		1-Naphthalenecarbonitrile	153	C11H7N	000086-53-3	87
5		1-Naphthalenecarbonitrile	153	C11H7N	000086-53-3	83



Data Path : Z:\svoasrv\HPCHEM1\BNA N\Data\BN022420\
Data File : BN009898.D
Acq On : 25 Feb 2020 13:17
Operator : CG/JU
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Misc :
ALS Vial : 36 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
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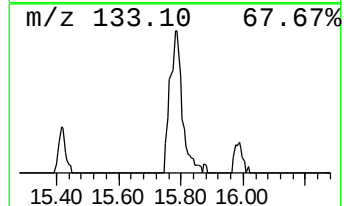
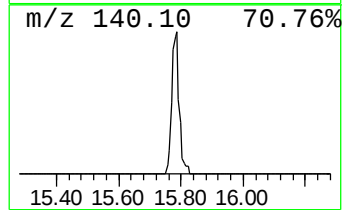
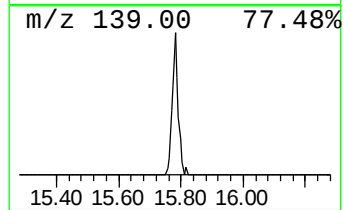
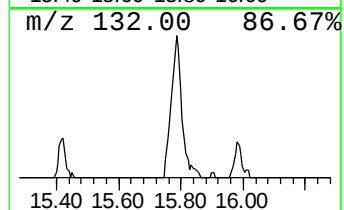
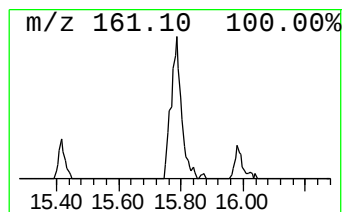
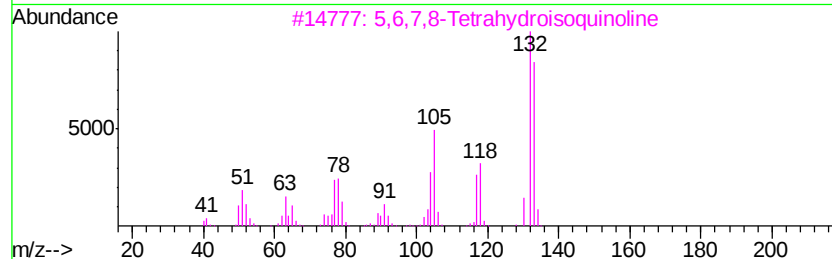
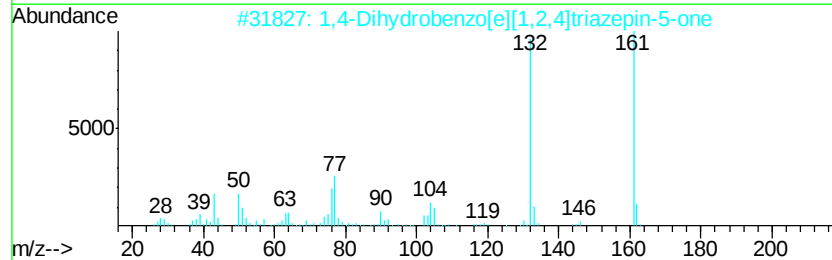
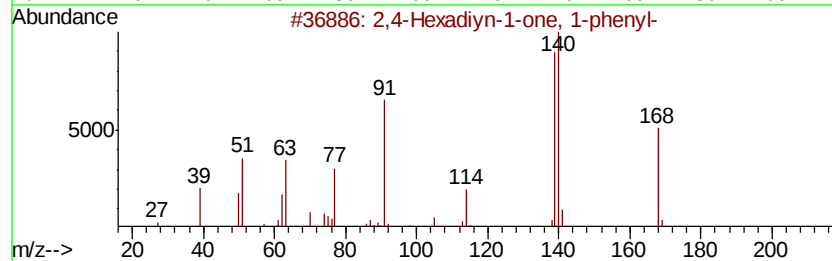
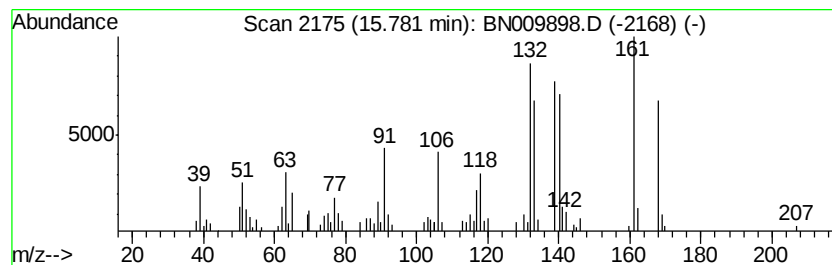
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 unknown-01 Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.78	3.66 ng/ul	236667	Phenanthrene-d10	16.79

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2,4-Hexadiyn-1-one, 1-phenyl-	168	C12H8O	000495-74-9	42
2			1,4-Dihydrobenzo[e][1,2,4]triazepin-5-one	161	C8H7N3O	057711-25-8	38
3			5,6,7,8-Tetrahydroisoquinoline	133	C9H11N	1000379-32-5	38
4			Quinoline, 1,2,3,4-tetrahydro-	133	C9H11N	000635-46-1	35
5			1(2H)-Acenaphthylenone	168	C12H8O	002235-15-6	25



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN022420\
Data File : BN009898.D
Acq On : 25 Feb 2020 13:17
Operator : CG/JU
Sample : L1582-07DL2 30X
Misc :
ALS Vial : 36 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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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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p-Cymene	7.52	2.1	ng/ul	62738	1	7.39	605749	20.0
Indane	7.75	8.6	ng/ul	261424	1	7.39	605749	20.0
Benzene, 1-propynyl-	7.91	23.8	ng/ul	719298	1	7.39	605749	20.0
Benzofuran, 2-met...	8.85	2.4	ng/ul	103950	2	10.13	871365	20.0
Bicyclo[2.2.1]hep...	9.55	3.5	ng/ul	151803	2	10.13	871365	20.0
Benzo[b]thiophene	10.30	8.7	ng/ul	379838	2	10.13	871365	20.0
Naphthalene, 1-me...	12.03	18.4	ng/ul	800344	2	10.13	871365	20.0
2-Naphthalenecarb...	14.18	2.2	ng/ul	119723	3	14.03	1073250	20.0
unknown-01	15.78	3.7	ng/ul	236667	4	16.79	1294530	20.0