

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN042920\
 Data File : BN010648.D
 Acq On : 29 Apr 2020 08:17
 Operator : CG/JU
 Sample : L2418-08DL 10X
 Misc :
 ALS Vial : 28 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 CB612DL

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-BN040920MA.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	2.876	11	15	25	rBV	85468	179890	2.42%	0.263%
2	2.958	25	29	35	rVB2	59589	87413	1.17%	0.128%
3	3.228	71	75	87	rVB	247139	393131	5.28%	0.575%
4	3.970	197	201	206	rBV	45607	63327	0.85%	0.093%
5	4.317	255	260	266	rBV	75793	125063	1.68%	0.183%
6	4.975	366	372	377	rBV5	16176	31078	0.42%	0.045%
7	5.293	421	426	433	rVB6	16949	34488	0.46%	0.050%
8	5.905	525	530	536	rVB	17147	29131	0.39%	0.043%
9	6.305	586	598	606	rBV2	85611	168158	2.26%	0.246%
10	6.564	636	642	644	rBV5	27948	40753	0.55%	0.060%
11	6.593	644	647	652	rVV2	48450	82991	1.11%	0.121%
12	6.781	672	679	685	rBV2	63532	120548	1.62%	0.176%
13	6.928	699	704	712	rBV	163412	259369	3.48%	0.379%
14	7.128	732	738	746	rBV	197144	332238	4.46%	0.486%
15	7.258	752	760	767	rVB8	12947	29526	0.40%	0.043%
16	7.581	806	815	823	rBV	1875444	2929223	39.35%	4.283%
17	7.652	823	827	831	rVV2	34002	53145	0.71%	0.078%
18	7.722	832	839	850	rVB	963808	1484600	19.94%	2.171%
19	7.816	850	855	860	rBV8	15527	29428	0.40%	0.043%
20	7.893	860	868	873	rVB7	36910	61918	0.83%	0.091%
21	7.946	873	877	884	rVB3	22333	41569	0.56%	0.061%
22	8.293	931	936	942	rVV	168683	260220	3.50%	0.380%
23	8.469	961	966	969	rBV4	28481	54407	0.73%	0.080%
24	8.552	974	980	991	rBV	3116440	4985010	66.97%	7.288%
25	8.675	996	1001	1004	rVV3	51214	88763	1.19%	0.130%
26	8.716	1004	1008	1018	rVV	195342	362459	4.87%	0.530%
27	8.811	1018	1024	1032	rVB	869211	1316780	17.69%	1.925%
28	8.928	1038	1044	1045	rBV4	20640	31085	0.42%	0.045%
29	8.958	1045	1049	1057	rVB8	48220	94133	1.26%	0.138%
30	9.081	1064	1070	1075	rBV5	43876	85402	1.15%	0.125%
31	9.193	1083	1089	1097	rBV	79841	129497	1.74%	0.189%
32	9.269	1097	1102	1107	rBV6	15796	33737	0.45%	0.049%
33	9.340	1110	1114	1118	rBV7	20302	29522	0.40%	0.043%
34	9.434	1124	1130	1136	rBV	166481	293529	3.94%	0.429%

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 Title : SVOA CALIBRATION

35	9.481	1136	1138	1143	rVB3	29747	42545	0.57%	0.062%
36	9.569	1146	1153	1156	rBV8	31279	68486	0.92%	0.100%
37	9.616	1156	1161	1168	rVB4	63422	125237	1.68%	0.183%
38	9.699	1168	1175	1176	rBV2	288421	413614	5.56%	0.605%
39	9.722	1176	1179	1182	rVV	511389	769799	10.34%	1.125%
40	9.763	1182	1186	1195	rVB	1470108	2277850	30.60%	3.330%
41	9.899	1201	1209	1215	rVB	126116	229084	3.08%	0.335%
42	9.969	1215	1221	1226	rBV	317460	571232	7.67%	0.835%
43	10.128	1243	1248	1257	rBV9	34271	84593	1.14%	0.124%
44	10.216	1257	1263	1272	rVV	606079	947446	12.73%	1.385%
45	10.340	1272	1284	1290	rVV	2678960	4418954	59.36%	6.461%
46	10.410	1290	1296	1301	rVV	1660188	2548226	34.23%	3.726%
47	10.475	1301	1307	1321	rVB	328471	687580	9.24%	1.005%
48	10.616	1321	1331	1335	rBV	31220	87268	1.17%	0.128%
49	10.693	1340	1344	1348	rVB7	22586	31904	0.43%	0.047%
50	10.916	1377	1382	1385	rBV5	24109	41187	0.55%	0.060%
51	10.981	1388	1393	1396	rVV6	20199	40817	0.55%	0.060%
52	11.028	1396	1401	1409	rVB10	29958	66941	0.90%	0.098%
53	11.175	1420	1426	1437	rVB	315850	558902	7.51%	0.817%
54	11.322	1446	1451	1457	rVB8	31154	59108	0.79%	0.086%
55	11.393	1459	1463	1468	rVB8	19207	30980	0.42%	0.045%
56	11.546	1485	1489	1495	rVB4	30288	51136	0.69%	0.075%
57	11.699	1506	1515	1519	rBV	78900	148035	1.99%	0.216%
58	11.846	1534	1540	1547	rVV	965898	1509289	20.27%	2.207%
59	11.910	1549	1551	1553	rVV3	27731	33120	0.44%	0.048%
60	11.952	1553	1558	1566	rVB3	150792	279631	3.76%	0.409%
61	12.087	1574	1581	1587	rVB	117692	202054	2.71%	0.295%
62	12.357	1622	1627	1631	rBV5	40961	62643	0.84%	0.092%
63	12.446	1638	1642	1647	rVB7	18506	32374	0.43%	0.047%
64	12.922	1717	1723	1729	rBV8	31406	63412	0.85%	0.093%
65	13.057	1743	1746	1754	rVB3	43051	65097	0.87%	0.095%
66	13.128	1754	1758	1762	rBV6	26625	46019	0.62%	0.067%
67	13.246	1774	1778	1785	rBV4	40017	68515	0.92%	0.100%
68	13.316	1785	1790	1797	rVB7	27365	53081	0.71%	0.078%
69	13.628	1838	1843	1847	rVV	535926	739010	9.93%	1.080%
70	13.669	1847	1850	1859	rVB	94034	142880	1.92%	0.209%
71	13.769	1861	1867	1870	rBV8	24375	50267	0.68%	0.073%

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72	13.893	1883	1888	1895	rBV	543957	870916	11.70%	1.273%
73	14.028	1907	1911	1915	rBV2	53022	79644	1.07%	0.116%
74	14.134	1924	1929	1934	rVB3	112628	157793	2.12%	0.231%
75	14.210	1935	1942	1950	rBV	4118096	6223908	83.61%	9.100%
76	14.451	1978	1983	1986	rBV7	26446	47872	0.64%	0.070%
77	14.910	2058	2061	2067	rVB5	26304	44155	0.59%	0.065%
78	14.969	2068	2071	2076	rVV	54275	82329	1.11%	0.120%
79	15.204	2106	2111	2117	rVB	774275	1083515	14.56%	1.584%
80	15.328	2128	2132	2137	rBV	71418	110262	1.48%	0.161%
81	15.463	2151	2155	2159	rBV7	50306	75355	1.01%	0.110%
82	15.728	2194	2200	2210	rVB	670650	971640	13.05%	1.421%
83	15.893	2223	2228	2233	rBV2	150156	247494	3.32%	0.362%
84	16.057	2252	2256	2261	rBV	38830	56476	0.76%	0.083%
85	16.234	2278	2286	2294	rBV	436809	675803	9.08%	0.988%
86	16.516	2330	2334	2340	rVB2	47136	64533	0.87%	0.094%
87	16.745	2370	2373	2375	rBV2	41592	51285	0.69%	0.075%
88	16.957	2402	2409	2420	rBV	5246568	7444195	100.00%	10.884%
89	17.051	2420	2425	2431	rBV2	841126	1202919	16.16%	1.759%
90	17.151	2438	2442	2448	rBV	63771	108900	1.46%	0.159%
91	17.881	2563	2566	2570	rBV6	29129	43395	0.58%	0.063%
92	18.157	2610	2613	2616	rBV5	43060	48673	0.65%	0.071%
93	18.645	2692	2696	2702	rVB3	121749	157700	2.12%	0.231%
94	19.357	2809	2817	2823	rBV	1012488	1405892	18.89%	2.055%
95	19.416	2823	2827	2835	rVV	189885	273090	3.67%	0.399%
96	21.163	3119	3124	3131	rBV	5720586	6850791	92.03%	10.016%
97	22.716	3385	3388	3394	rVB	395230	468599	6.29%	0.685%
98	23.192	3465	3469	3474	rVB	501287	815121	10.95%	1.192%
99	23.257	3477	3480	3485	rVV	485229	710961	9.55%	1.039%
100	23.327	3486	3492	3501	rVB	3321625	6034401	81.06%	8.823%

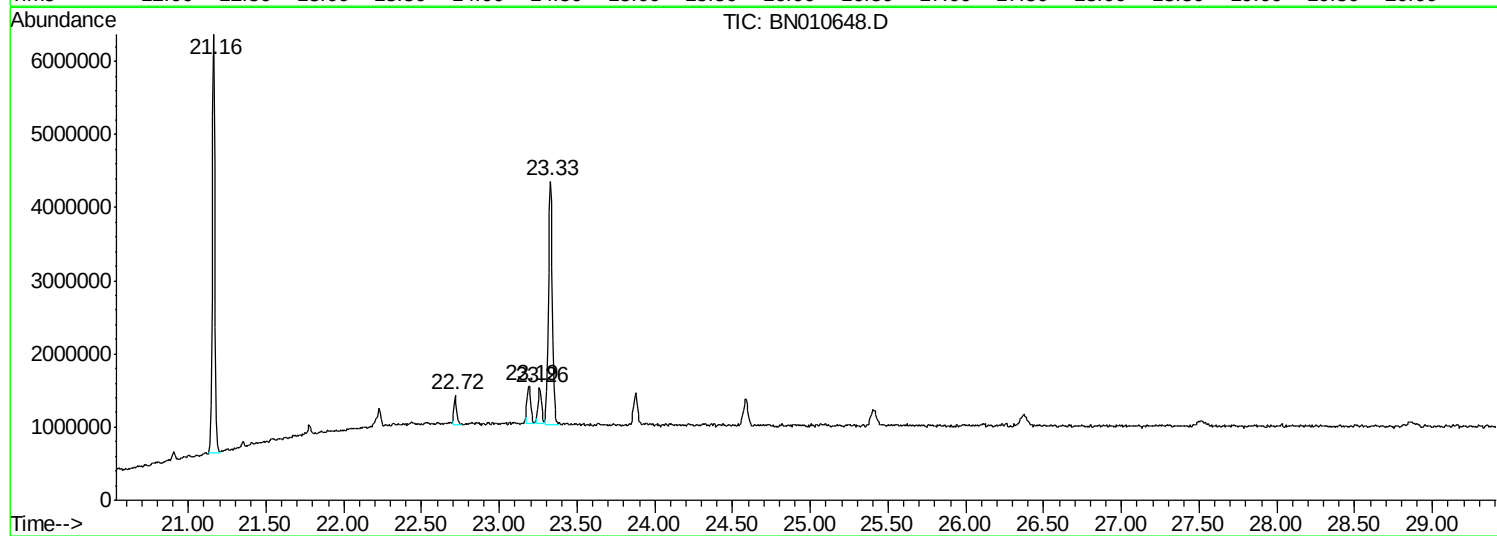
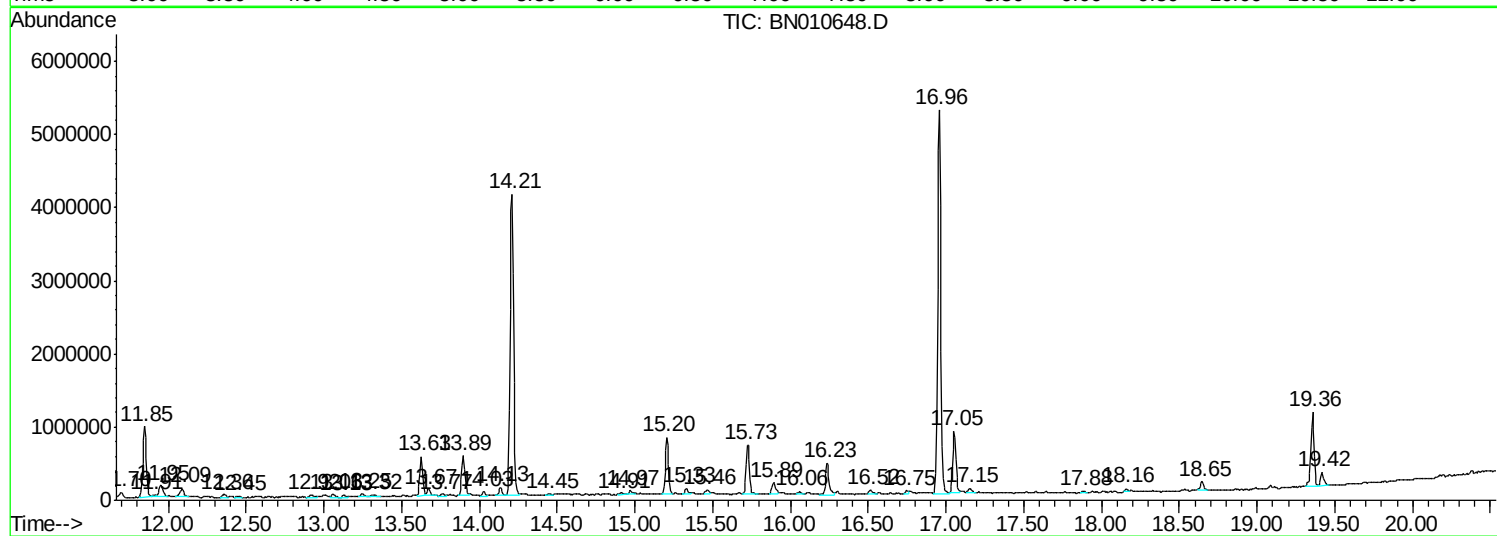
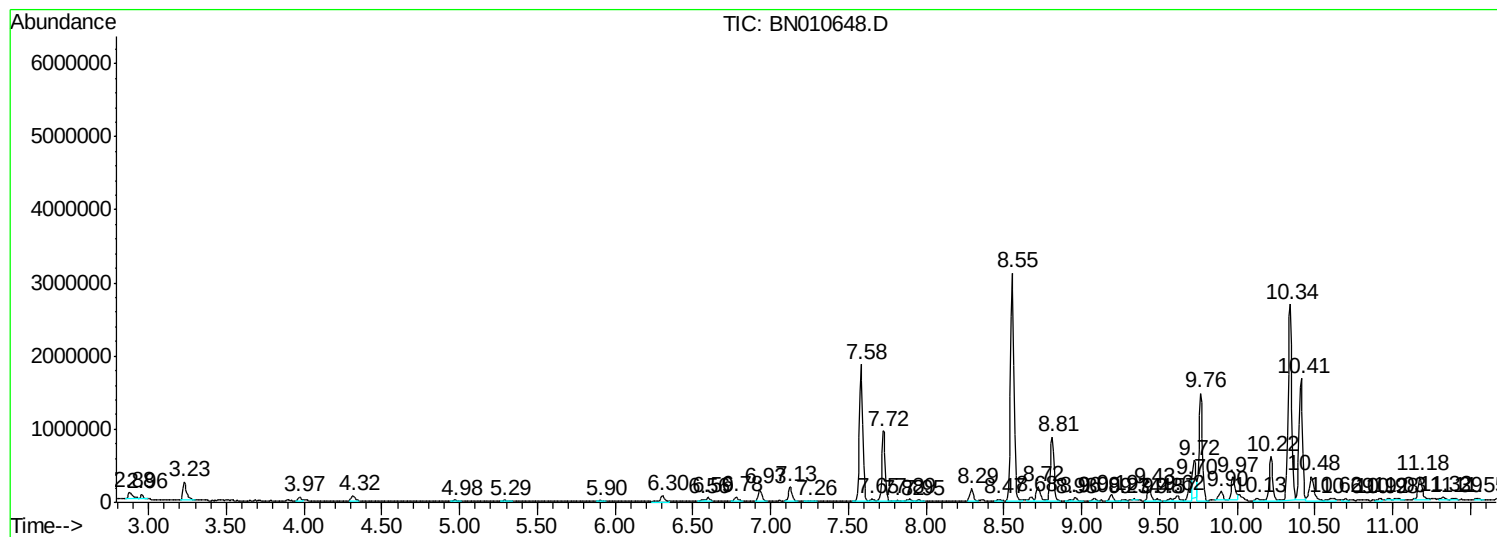
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Quant Title :

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



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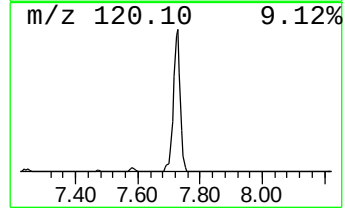
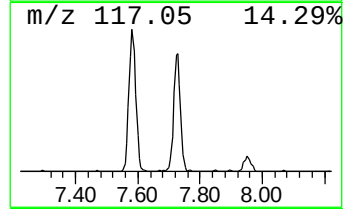
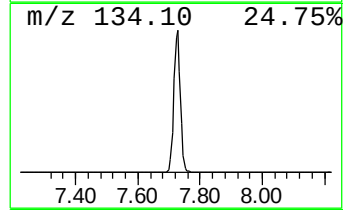
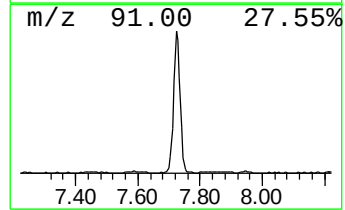
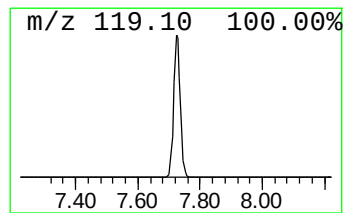
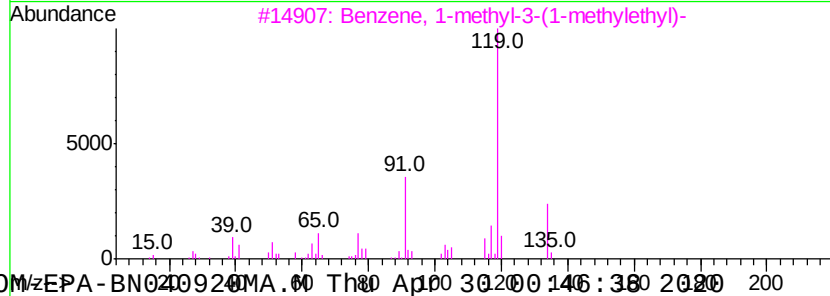
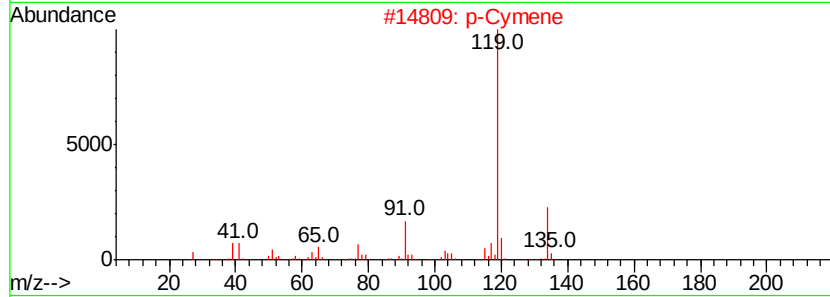
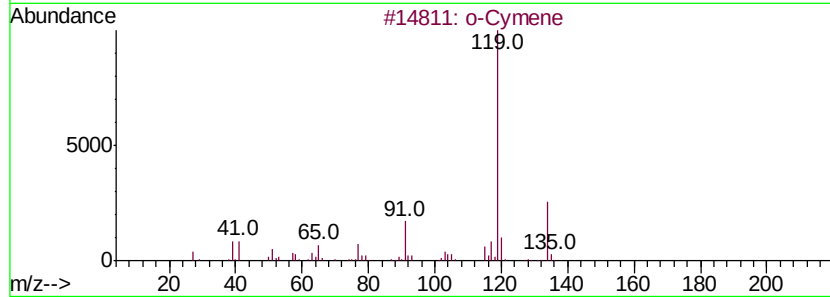
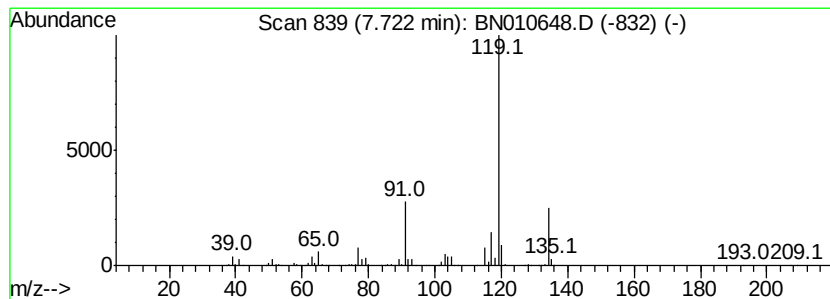
Quant Title :

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

Peak Number 1 o-Cymene Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.72	10.14 ng/ul	1484600	1,4-Dichlorobenzene-d4	7.58

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		o-Cymene	134	C10H14	000527-84-4	94
2		p-Cymene	134	C10H14	000099-87-6	94
3		Benzene, 1-methyl-3-(1-methylethyl-)	134	C10H14	000535-77-3	94
4		Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	91
5		o-Cymene	134	C10H14	000527-84-4	91



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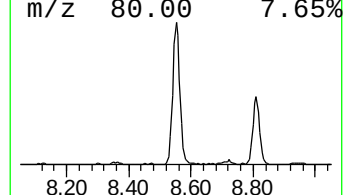
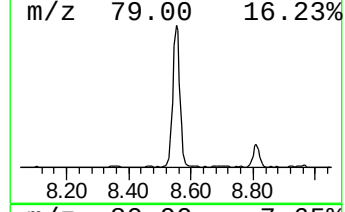
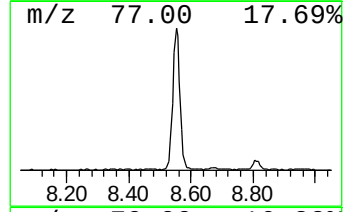
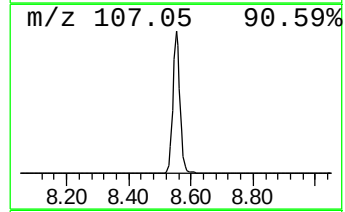
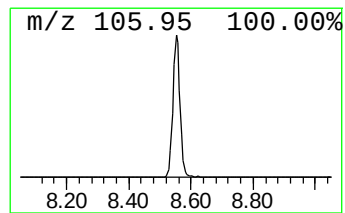
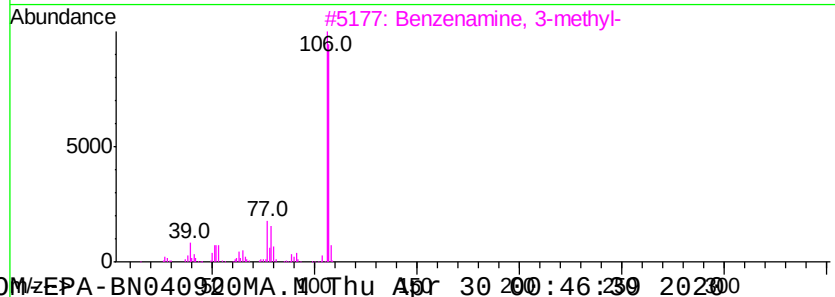
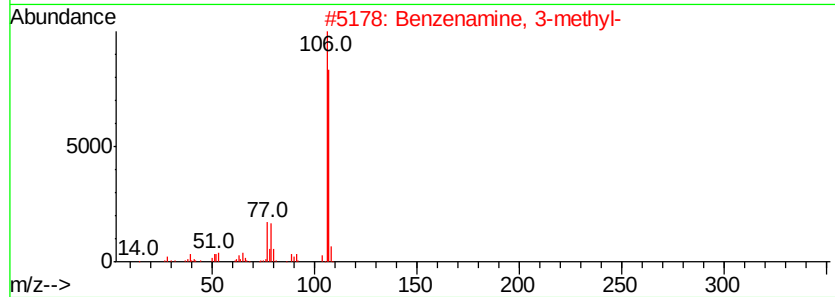
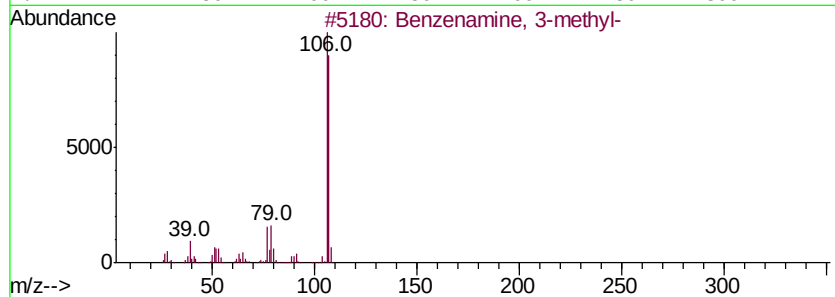
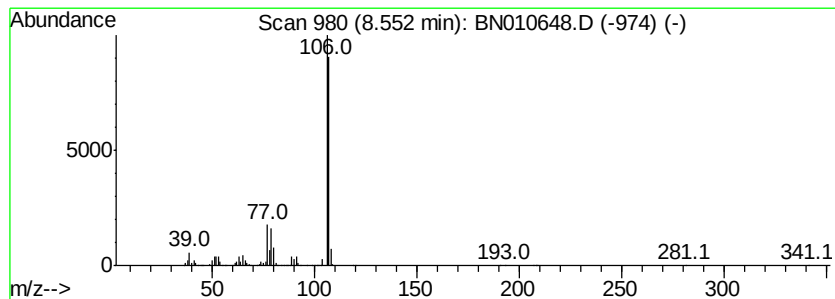
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 Benzenamine, 3-methyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.	
8.55	34.04 ng/ul	4985010	1,4-Dichlorobenzene-d4	7.58	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzenamine, 3-methyl-	107	C7H9N	000108-44-1	97
2	Benzenamine, 3-methyl-	107	C7H9N	000108-44-1	97
3	Benzenamine, 3-methyl-	107	C7H9N	000108-44-1	97
4	Benzenamine, 3-methyl-	107	C7H9N	000108-44-1	94
5	o-Toluidine	107	C7H9N	000095-53-4	94



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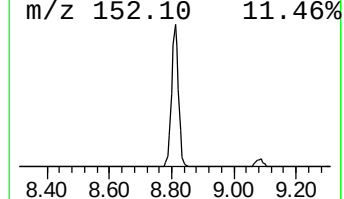
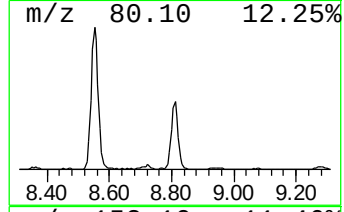
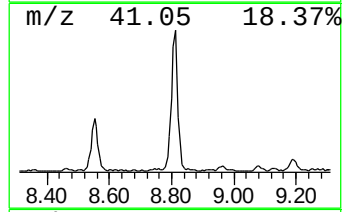
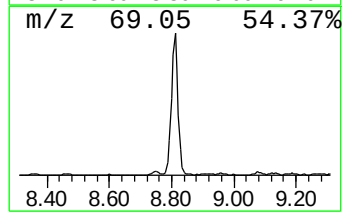
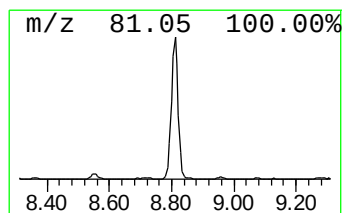
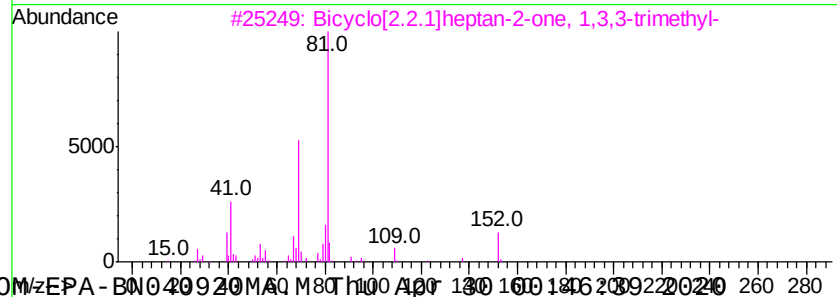
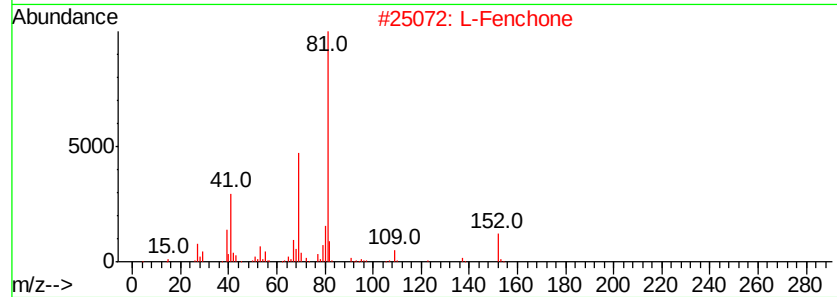
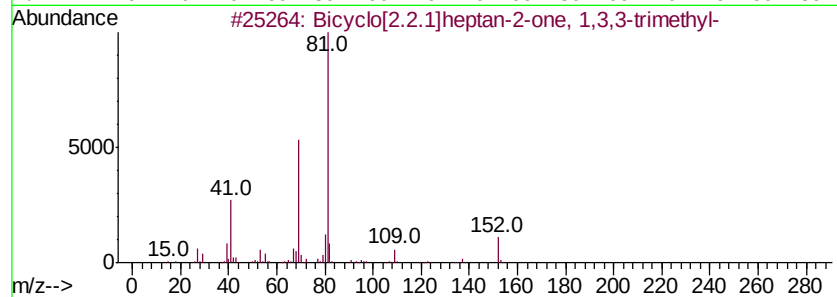
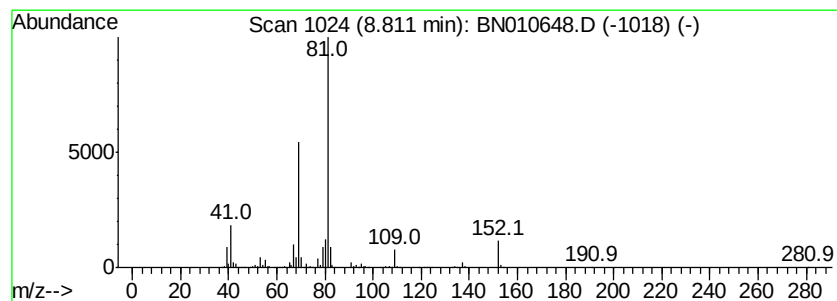
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 Bicyclo[2.2.1]heptan-2-one, ... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.81	8.99 ng/ul	1316780	1,4-Dichlorobenzene-d4	7.58

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Bicyclo[2.2.1]heptan-2-one, 1,3,...	152	C10H16O	001195-79-5	95
2	L-Fenchone	152	C10H16O	007787-20-4	94
3	Bicyclo[2.2.1]heptan-2-one, 1,3,...	152	C10H16O	001195-79-5	94
4	D-Fenchone	152	C10H16O	004695-62-9	91
5	L-Fenchone	152	C10H16O	007787-20-4	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN042920\
Data File : BN010648.D
Acq On : 29 Apr 2020 08:17
Operator : CG/JU
Sample : L2418-08DL 10X
Misc :
ALS Vial : 28 Sample Multiplier: 1

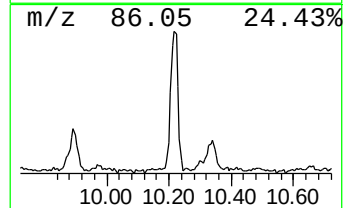
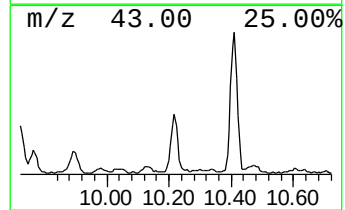
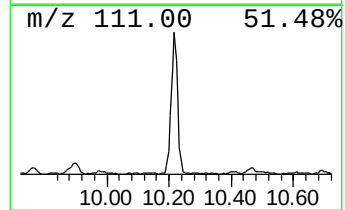
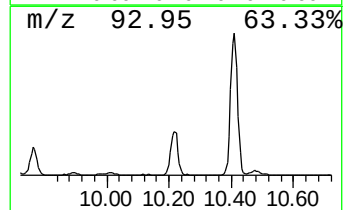
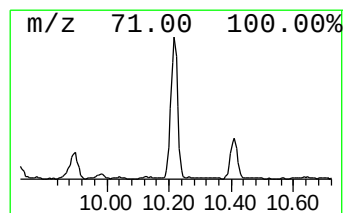
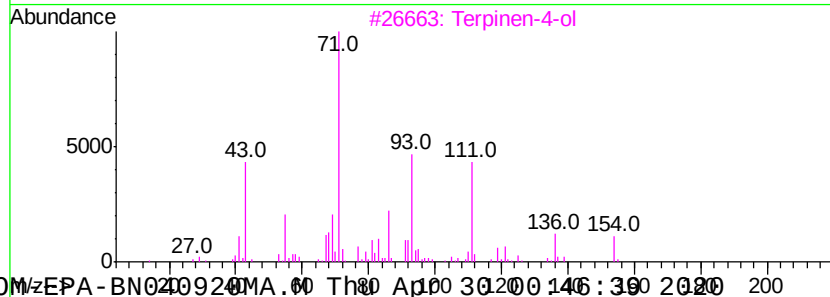
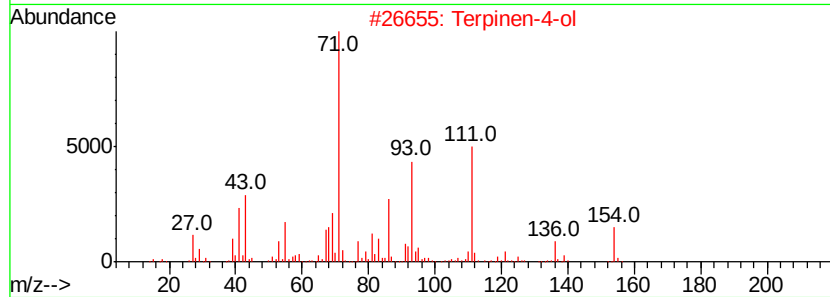
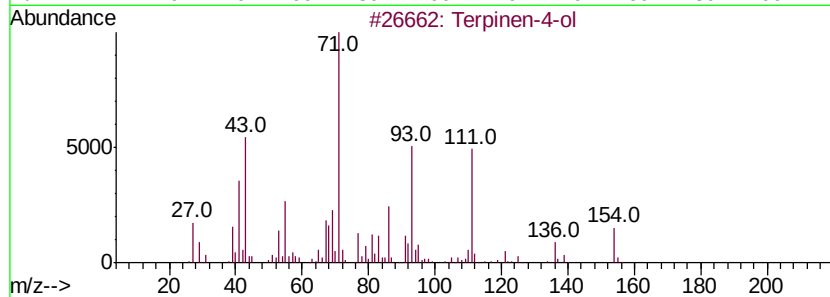
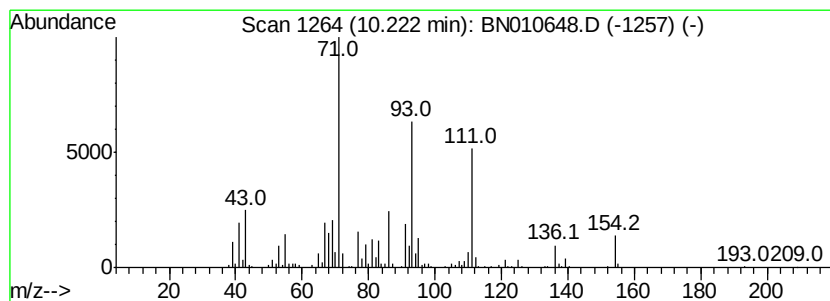
Instrument :
BNA_N
ClientSampleId :
CB612DL

Quant Title :

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

Peak Number 4 Terpinen-4-ol Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.22	4.29 ng/ul	947446	Napthalene-d8	10.34
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Terpinen-4-ol	154 C10H18O	000562-74-3	93
2	Terpinen-4-ol	154 C10H18O	000562-74-3	90
3	Terpinen-4-ol	154 C10H18O	000562-74-3	90
4	3-Cyclohexen-1-ol, 4-methyl-1-(1...	154 C10H18O	020126-76-5	81
5	Terpinen-4-ol	154 C10H18O	000562-74-3	74



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN042920\
Data File : BN010648.D
Acq On : 29 Apr 2020 08:17
Operator : CG/JU
Sample : L2418-08DL 10X
Misc :
ALS Vial : 28 Sample Multiplier: 1

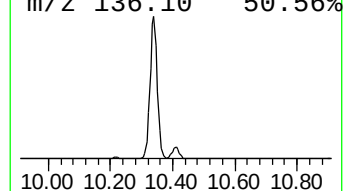
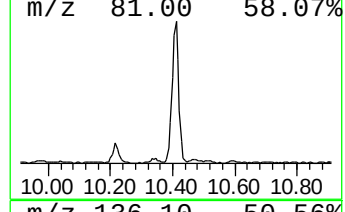
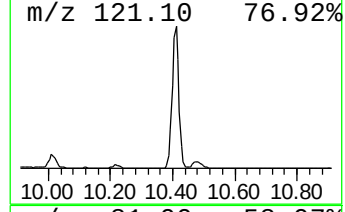
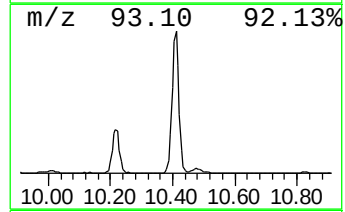
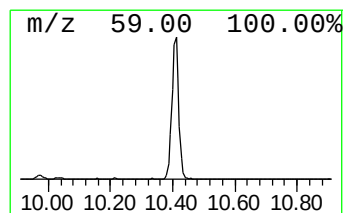
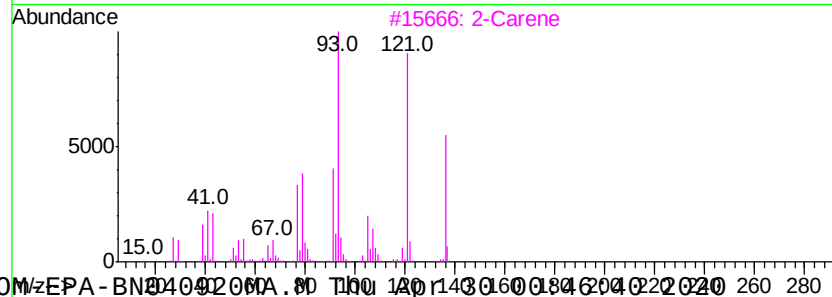
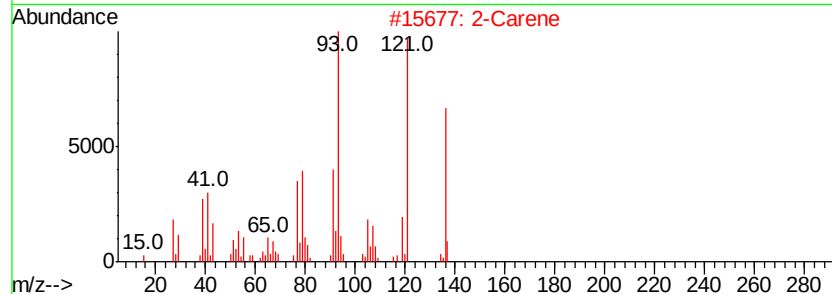
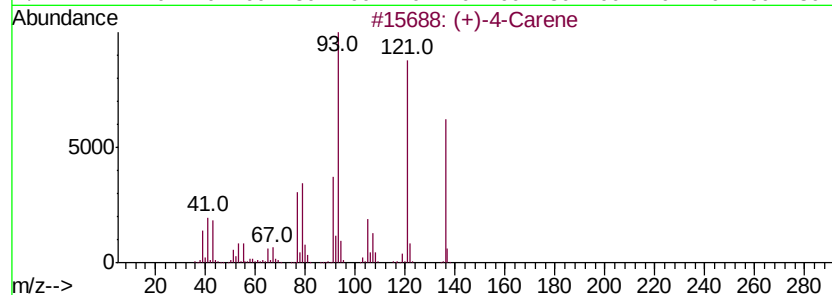
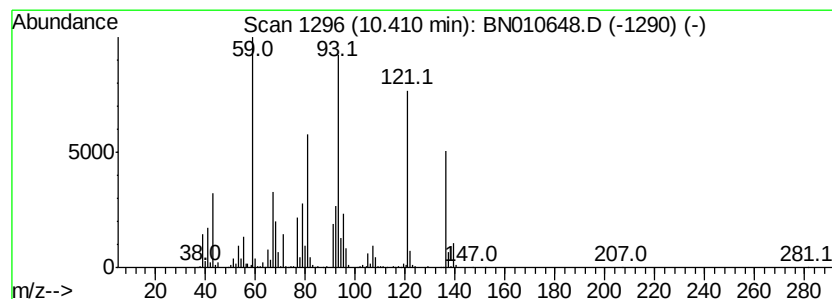
Instrument :
BNA_N
ClientSampleId :
CB612DL

Quant Title :

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

Peak Number 5 (+)-4-Carene Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.	
10.41	11.53 ng/ul	2548230	Napthalene-d8	10.34	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	(+)-4-Carene	136	C10H16	029050-33-7	83
2	2-Carene	136	C10H16	000554-61-0	83
3	2-Carene	136	C10H16	000554-61-0	78
4	L-.alpha.-Terpineol	154	C10H18O	010482-56-1	72
5	(1R)-2,6,6-Trimethylbicyclo[3.1....	136	C10H16	007785-70-8	70



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN042920\
Data File : BN010648.D
Acq On : 29 Apr 2020 08:17
Operator : CG/JU
Sample : L2418-08DL 10X
Misc :
ALS Vial : 28 Sample Multiplier: 1

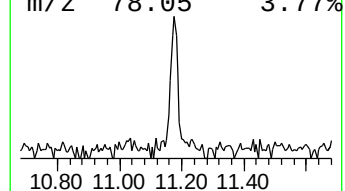
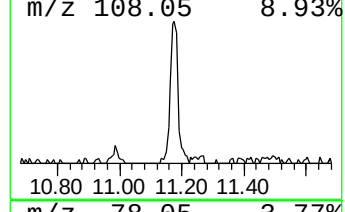
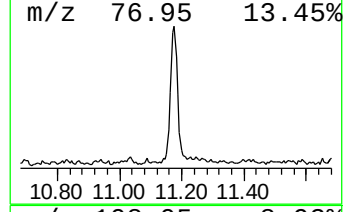
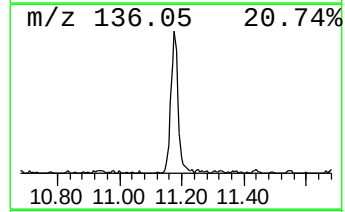
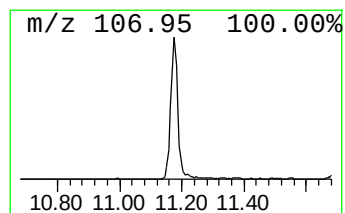
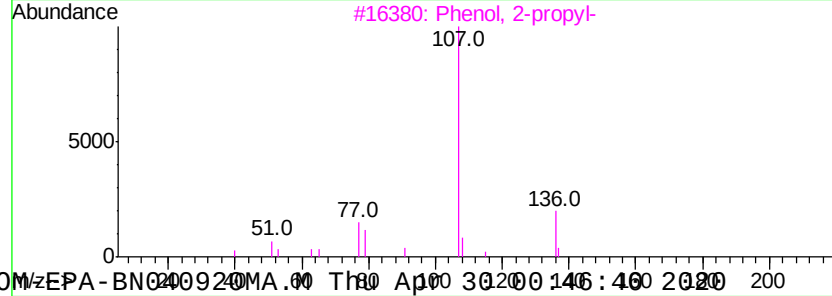
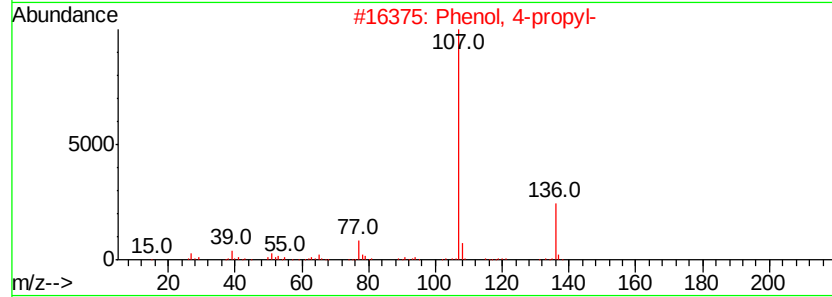
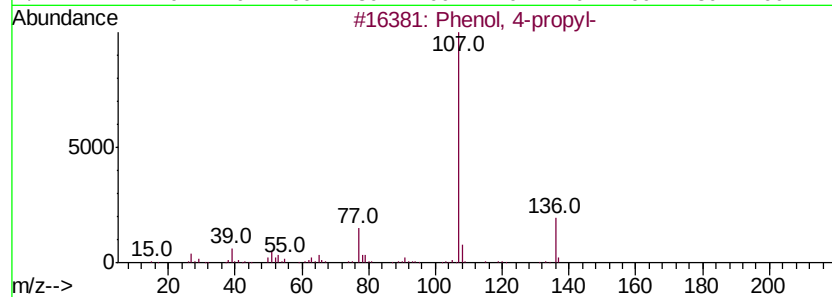
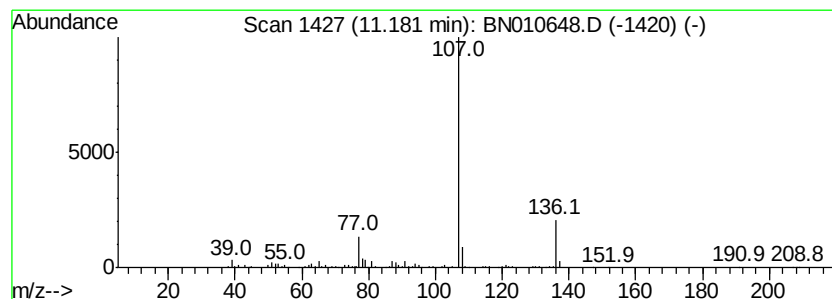
Instrument :
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ClientSampleId :
CB612DL

Quant Title :

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

Peak Number 6 Phenol, 4-propyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.	
11.18	2.53 ng/ul	558902	Napthalene-d8	10.34	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, 4-propyl-	136	C9H12O	000645-56-7	87
2	Phenol, 4-propyl-	136	C9H12O	000645-56-7	87
3	Phenol, 2-propyl-	136	C9H12O	000644-35-9	72
4	Furan, 2-(1-pentenyl)-, (Z)-	136	C9H12O	020992-60-3	56
5	Pyridine, 2,5-dimethyl-	107	C7H9N	000589-93-5	53



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN042920\
Data File : BN010648.D
Acq On : 29 Apr 2020 08:17
Operator : CG/JU
Sample : L2418-08DL 10X
Misc :
ALS Vial : 28 Sample Multiplier: 1

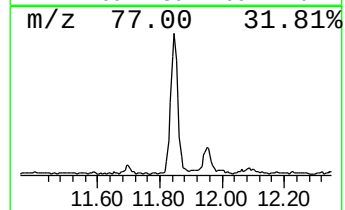
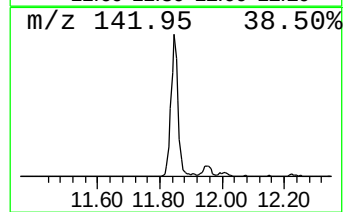
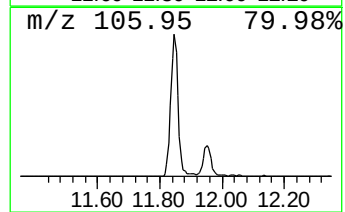
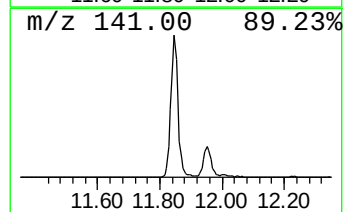
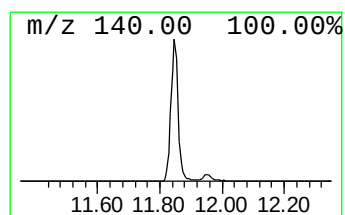
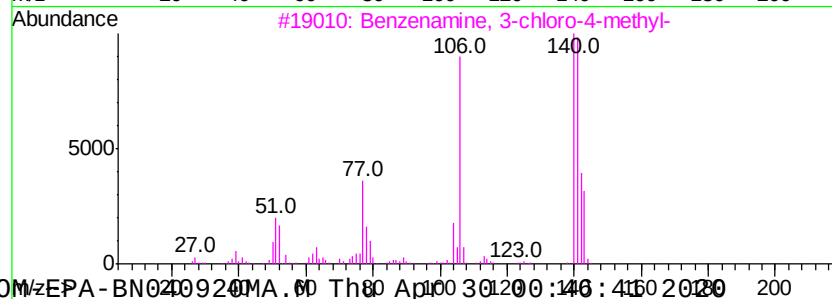
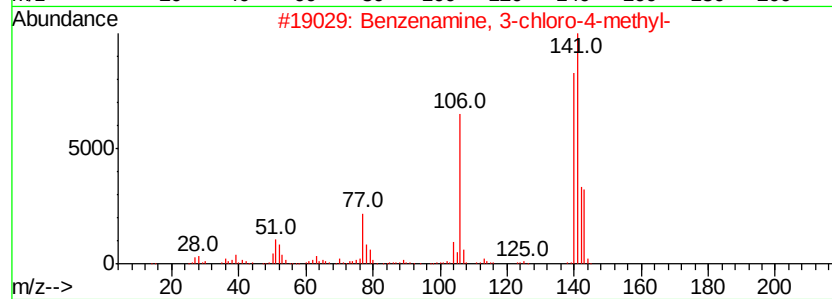
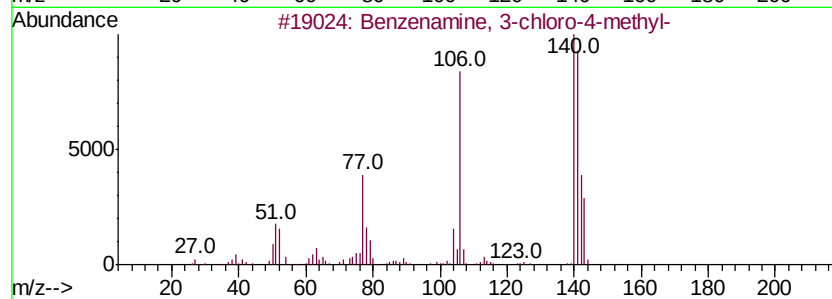
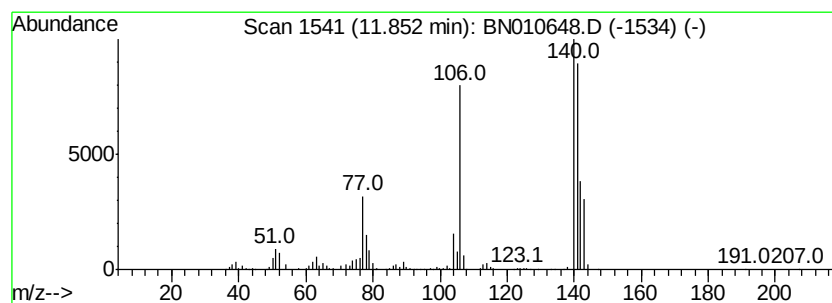
Instrument :
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ClientSampleId :
CB612DL

Quant Title :

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

Peak Number 7 Benzenamine, 3-chloro-4-met... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.	
11.85	6.83 ng/ul	1509290	Napthalene-d8	10.34	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzenamine, 3-chloro-4-methyl-	141	C7H8ClN	000095-74-9	97
2	Benzenamine, 3-chloro-4-methyl-	141	C7H8ClN	000095-74-9	95
3	Benzenamine, 3-chloro-4-methyl-	141	C7H8ClN	000095-74-9	94
4	Benzenamine, 2-chloro-4-methyl-	141	C7H8ClN	000615-65-6	91
5	Benzenamine, 5-chloro-2-methyl-	141	C7H8ClN	000095-79-4	74



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN042920\
Data File : BN010648.D
Acq On : 29 Apr 2020 08:17
Operator : CG/JU
Sample : L2418-08DL 10X
Misc :
ALS Vial : 28 Sample Multiplier: 1

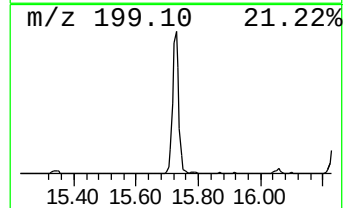
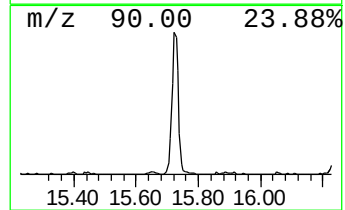
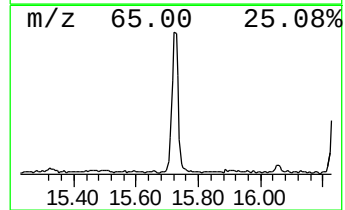
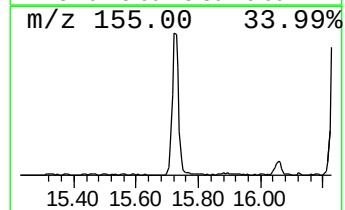
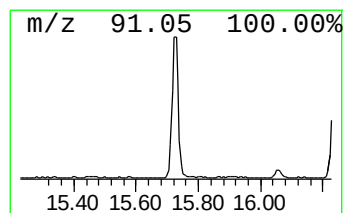
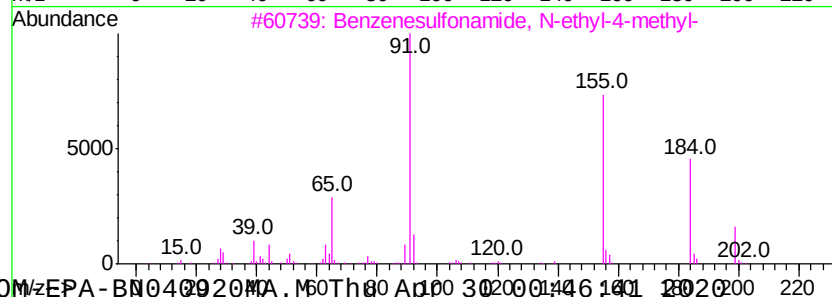
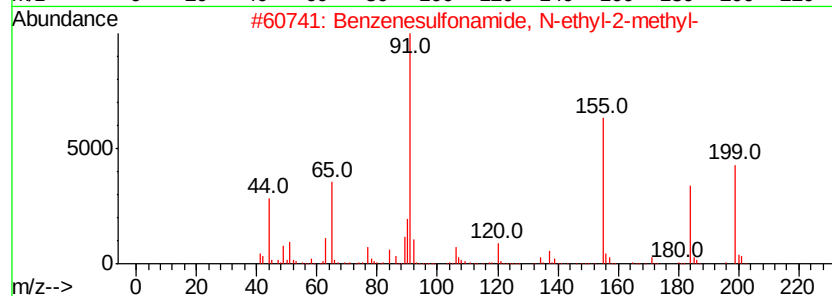
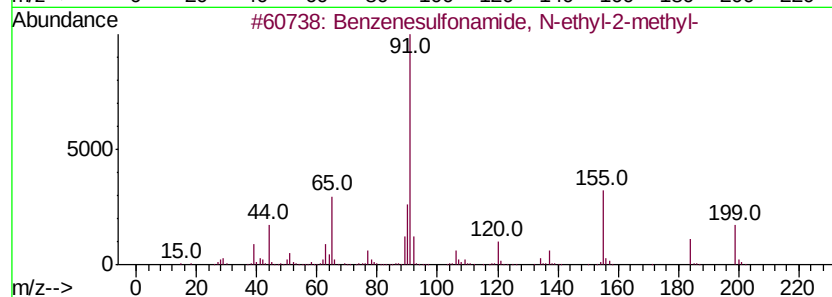
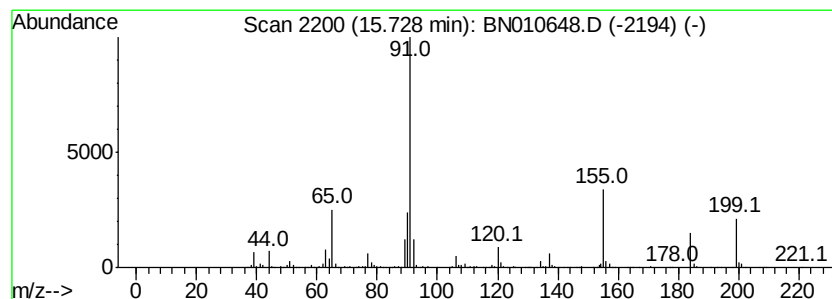
Instrument :
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ClientSampleId :
CB612DL

Quant Title :

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

Peak Number 8 Benzenesulfonamide, N-ethyl... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.	
15.73	2.61 ng/ul	971640	Phenanthrene-d10	16.96	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzenesulfonamide, N-ethyl-2-me...	199	C9H13NO2S	001077-56-1	95
2	Benzenesulfonamide, N-ethyl-2-me...	199	C9H13NO2S	001077-56-1	50
3	Benzenesulfonamide, N-ethyl-4-me...	199	C9H13NO2S	000080-39-7	50
4	N3, N.alpha.-bis*ptoluenesulfony...	462	C20H22N4O5S2	1000130-21-9	47
5	(O-p-Tosylisonitroso)malononitrile	249	C10H7N3O3S	020893-01-0	43



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN042920\
Data File : BN010648.D
Acq On : 29 Apr 2020 08:17
Operator : CG/JU
Sample : L2418-08DL 10X
Misc :
ALS Vial : 28 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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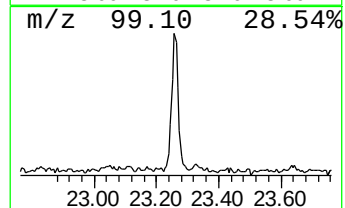
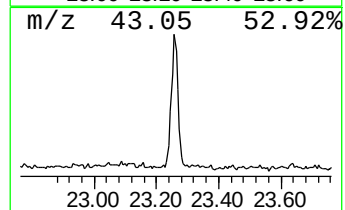
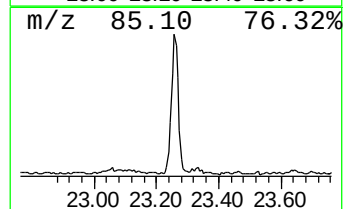
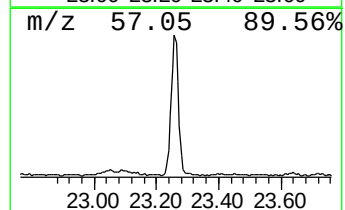
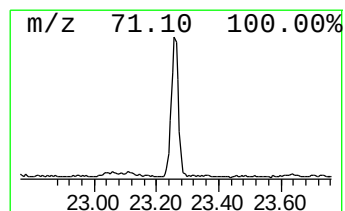
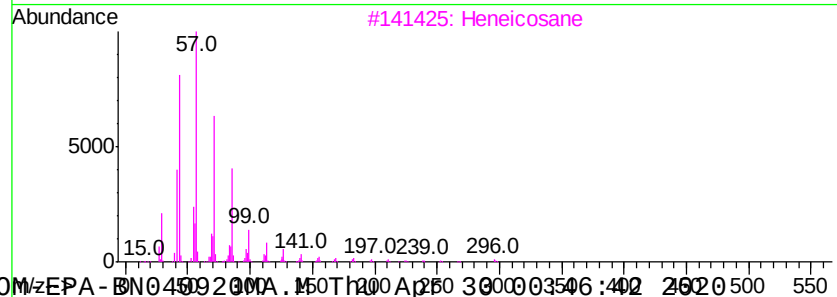
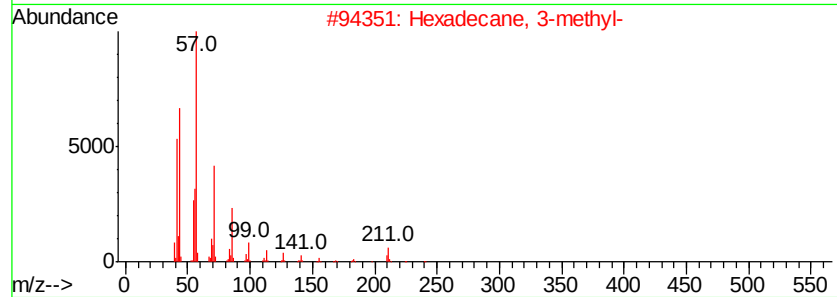
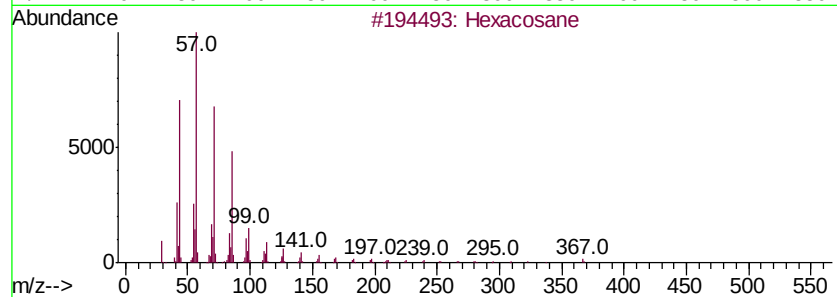
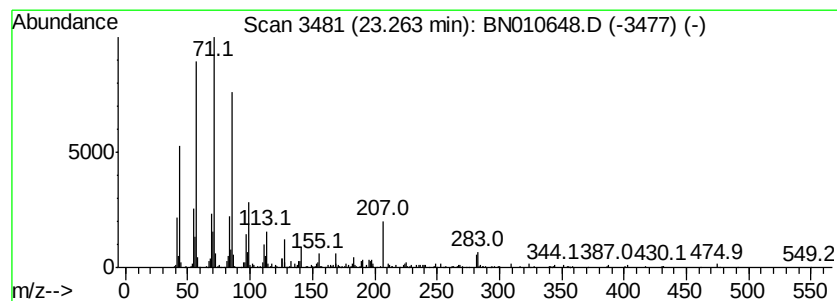
Quant Title :

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

Peak Number 9 (DEL) Alkane: Straight-Chai... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.26	2.36 ng/ul	710961	Perylene-d12	23.33

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexacosane	366	C26H54	000630-01-3	87
2	Hexadecane, 3-methyl-	240	C17H36	006418-43-5	87
3	Heneicosane	296	C21H44	000629-94-7	83
4	Eicosane, 10-methyl-	296	C21H44	054833-23-7	83
5	Tetracosane	338	C24H50	000646-31-1	83



Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN042920\
Data File : BN010648.D
Acq On : 29 Apr 2020 08:17
Operator : CG/JU
Sample : L2418-08DL 10X
Misc :
ALS Vial : 28 Sample Multiplier: 1

Instrument :
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ClientSampleId :
CB612DL

Quant Title :

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
o-Cymene	7.72	10.1	ng/ul	1484600	1	7.58	2929220	20.0
Benzenamine, 3-me...	8.55	34.0	ng/ul	4985010	1	7.58	2929220	20.0
Bicyclo[2.2.1]hep...	8.81	9.0	ng/ul	1316780	1	7.58	2929220	20.0
Terpinen-4-ol	10.22	4.3	ng/ul	947446	2	10.34	4418950	20.0
(+)-4-Carene	10.41	11.5	ng/ul	2548230	2	10.34	4418950	20.0
Phenol, 4-propyl-	11.18	2.5	ng/ul	558902	2	10.34	4418950	20.0
Benzenamine, 3-ch...	11.85	6.8	ng/ul	1509290	2	10.34	4418950	20.0
Benzenesulfonamid...	15.73	2.6	ng/ul	971640	4	16.96	7444200	20.0
(DEL) Alkane: Str...	23.26	2.4	ng/ul	710961	6	23.33	6034400	20.0