LSC Area Percent Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM060120\
Data File : BM026197.D
Acq On : 01 Jun 2020 13:16
Operator : JU/CG
Sample : L2820-08DL 5X
Misc :
ALS Vial : 3 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : CB638DL

Integration Parameters: LSCINT.P
Integrator: RTE
Smoothing : OFF Filtering: 5
Sampling : 1 Min Area: 0.1 % of largest Peak
Start Thrs: 0.2 Max Peaks: 100
Stop Thrs : 0 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
Peak separation: 5

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

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.057	27	29	31	rBV2	10297	9861	0.49%	0.052%
2	3.087	31	34	43	rVB	105531	129491	6.43%	0.678%
3	3.599	118	121	124	rVB4	4527	6179	0.31%	0.032%
4	3.751	145	147	150	rVV2	8631	9837	0.49%	0.052%
5	3.840	159	162	166	rBV	21928	23580	1.17%	0.123%
6	4.181	216	220	226	rVB	35288	50982	2.53%	0.267%
7	4.469	267	269	273	rVB5	5480	6935	0.34%	0.036%
8	4.834	329	331	337	rVB5	6503	9262	0.46%	0.048%
9	5.528	446	449	454	rVB5	5042	7985	0.40%	0.042%
10	5.798	487	495	499	rBV3	5032	10022	0.50%	0.052%
11	6.181	552	560	564	rBV6	13297	22166	1.10%	0.116%
12	6.475	606	610	616	rVV3	22302	38486	1.91%	0.202%
13	6.675	639	644	650	rBV2	23770	39240	1.95%	0.205%
14	6.828	665	670	682	rBV	71192	118163	5.87%	0.619%
15	7.016	696	702	710	rVB	81813	136100	6.76%	0.713%
16	7.475	773	780	787	rVV	563081	884035	43.92%	4.629%
17	7.551	787	793	797	rVV6	9678	18997	0.94%	0.099%
18	7.616	797	804	813	rVB	455784	668810	33.23%	3.502%
19	7.698	813	818	823	rBV7	3845	7415	0.37%	0.039%
20	7.781	828	832	837	rVV4	12513	17717	0.88%	0.093%
21	7.845	839	843	847	rVB4	6545	8941	0.44%	0.047%
22	8.192	896	902	908	rVV	65102	104984	5.22%	0.550%
23	8.251	908	912	915	rBV5	4401	6221	0.31%	0.033%
24	8.357	926	930	935	rBV6	6065	13037	0.65%	0.068%
25	8.451	940	946	961	rVV	873619	1356908	67.42%	7.105%
26	8.569	963	966	969	rVV2	13125	20642	1.03%	0.108%
27	8.616	969	974	982	rVV	83894	153954	7.65%	0.806%
28	8.704	982	989	995	rVV	245725	391840	19.47%	2.052%
29	8.816	1003	1008	1011	rVV7	6020	11384	0.57%	0.060%
30	8.851	1011	1014	1021	rVB8	8659	15005	0.75%	0.079%
31	8.975	1029	1035	1038	rBV7	10399	18115	0.90%	0.095%
32	9.086	1048	1054	1062	rBV3	18519	34556	1.72%	0.181%
33	9.233	1075	1079	1083	rBV4	6206	8891	0.44%	0.047%
34	9.263	1083	1084	1090	rVB6	4969	7228	0.36%	0.038%

LSC Area Percent Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM060120\
Data File : BM026197.D
Acq On : 01 Jun 2020 13:16
Operator : JU/CG
Sample : L2820-08DL 5X
Misc :
ALS Vial : 3 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : CB638DL

Integration Parameters: LSCINT.P
Integrator: RTE
Smoothing : OFF Filtering: 5
Sampling : 1 Min Area: 0.1 % of largest Peak
Start Thrs: 0.2 Max Peaks: 100
Stop Thrs : 0 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

35	9.333	1090	1096	1100	rBV2	57919	98173	4.88%	0.514%
36	9.463	1111	1118	1120	rBV6	8305	14261	0.71%	0.075%
37	9.510	1121	1126	1134	rVV6	17408	38704	1.92%	0.203%
38	9.622	1134	1145	1148	rVV	606274	1057708	52.55%	5.538%
39	9.657	1148	1151	1159	rVV	433342	648572	32.23%	3.396%
40	9.733	1159	1164	1166	rVV6	4549	6993	0.35%	0.037%
41	9.786	1166	1173	1180	rVB3	41210	78576	3.90%	0.411%
42	9.869	1180	1187	1193	rBV	124095	227697	11.31%	1.192%
43	9.916	1193	1195	1205	rVB4	26418	47434	2.36%	0.248%
44	10.033	1207	1215	1220	rBV8	9017	22679	1.13%	0.119%
45	10.116	1223	1229	1235	rBV	73620	119221	5.92%	0.624%
46	10.233	1242	1249	1256	rBV	745088	1158506	57.56%	6.066%
47	10.304	1256	1261	1266	rVB6	23656	43971	2.18%	0.230%
48	10.380	1266	1274	1286	rBV	116137	236344	11.74%	1.238%
49	10.510	1289	1296	1302	rBV	6860	18699	0.93%	0.098%
50	10.592	1308	1310	1315	rVB6	4335	6450	0.32%	0.034%
51	10.816	1344	1348	1354	rBV7	5651	12538	0.62%	0.066%
52	10.939	1366	1369	1376	rVB9	5704	9271	0.46%	0.049%
53	11.080	1386	1393	1405	rVV	80699	149150	7.41%	0.781%
54	11.216	1411	1416	1417	rBV5	4466	6290	0.31%	0.033%
55	11.439	1450	1454	1460	rVB7	7476	14338	0.71%	0.075%
56	11.598	1474	1481	1487	rBV	17925	31729	1.58%	0.166%
57	11.751	1500	1507	1514	rBV	281501	446639	22.19%	2.339%
58	11.857	1521	1525	1532	rVB2	43127	72313	3.59%	0.379%
59	11.992	1540	1548	1553	rBV4	15716	30583	1.52%	0.160%
60	12.263	1590	1594	1598	rBV6	8019	13104	0.65%	0.069%
61	12.474	1623	1630	1633	rBV6	6624	12750	0.63%	0.067%
62	12.827	1684	1690	1698	rBV7	9646	21672	1.08%	0.113%
63	12.910	1700	1704	1708	rVV6	7903	14222	0.71%	0.074%
64	12.963	1711	1713	1719	rVB5	7697	12143	0.60%	0.064%
65	13.163	1742	1747	1750	rBV6	8069	12606	0.63%	0.066%
66	13.486	1800	1802	1806	rBV5	5427	6982	0.35%	0.037%
67	13.545	1806	1812	1816	rVV	188179	270915	13.46%	1.419%
68	13.586	1816	1819	1825	rVB	33433	50402	2.50%	0.264%
69	13.680	1831	1835	1840	rVB7	6213	6994	0.35%	0.037%
70	13.727	1840	1843	1848	rVB4	7915	12717	0.63%	0.067%
71	13.804	1850	1856	1863	rBV	194039	287585	14.29%	1.506%

LSC Area Percent Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM060120\
Data File : BM026197.D
Acq On : 01 Jun 2020 13:16
Operator : JU/CG
Sample : L2820-08DL 5X
Misc :
ALS Vial : 3 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : CB638DL

Integration Parameters: LSCINT.P
Integrator: RTE
Smoothing : OFF Filtering: 5
Sampling : 1 Min Area: 0.1 % of largest Peak
Start Thrs: 0.2 Max Peaks: 100
Stop Thrs : 0 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

72	13.945	1875	1880	1883	rBV6	4794	7943	0.39%	0.042%
73	14.045	1892	1897	1902	rBV6	24211	33674	1.67%	0.176%
74	14.115	1903	1909	1919	rBV	963719	1450268	72.06%	7.594%
75	14.368	1946	1952	1956	rBV6	11228	20024	0.99%	0.105%
76	14.721	2010	2012	2020	rBV9	3940	7519	0.37%	0.039%
77	14.827	2024	2030	2035	rBV3	16268	28331	1.41%	0.148%
78	14.886	2035	2040	2044	rBV	12912	19277	0.96%	0.101%
79	15.121	2074	2080	2086	rBV	258052	374125	18.59%	1.959%
80	15.257	2098	2103	2111	rBV3	43806	85058	4.23%	0.445%
81	15.380	2117	2124	2128	rBV9	16739	32205	1.60%	0.169%
82	15.592	2154	2160	2162	rBV6	6216	10739	0.53%	0.056%
83	15.645	2163	2169	2179	rVV	165016	243835	12.12%	1.277%
84	15.739	2183	2185	2188	rVB4	5852	6352	0.32%	0.033%
85	15.815	2193	2198	2210	rVV2	35405	71465	3.55%	0.374%
86	15.980	2221	2226	2230	rVV2	9400	16020	0.80%	0.084%
87	16.033	2230	2235	2237	rVV6	6756	11997	0.60%	0.063%
88	16.157	2250	2256	2265	rBV	94777	142596	7.09%	0.747%
89	16.433	2299	2303	2306	rBV5	10489	15995	0.79%	0.084%
90	16.462	2306	2308	2317	rVB7	16505	27318	1.36%	0.143%
91	16.656	2338	2341	2342	rBV2	8544	7772	0.39%	0.041%
92	16.868	2371	2377	2388	rBV	1274610	1743722	86.64%	9.130%
93	16.968	2388	2394	2402	rBV	280720	409951	20.37%	2.147%
94	17.074	2407	2412	2419	rVV	42380	65515	3.26%	0.343%
95	18.068	2578	2581	2587	rBV8	9985	16455	0.82%	0.086%
96	18.556	2661	2664	2668	rVB4	21946	24541	1.22%	0.129%
97	19.268	2780	2785	2791	rBV	341181	465604	23.13%	2.438%
98	19.333	2793	2796	2803	rVB2	44716	66370	3.30%	0.348%
99	21.068	3086	3091	3098	rBV	1577672	1994805	99.12%	10.445%
100	23.185	3445	3451	3459	rVB2	1150086	2012615	100.00%	10.538%

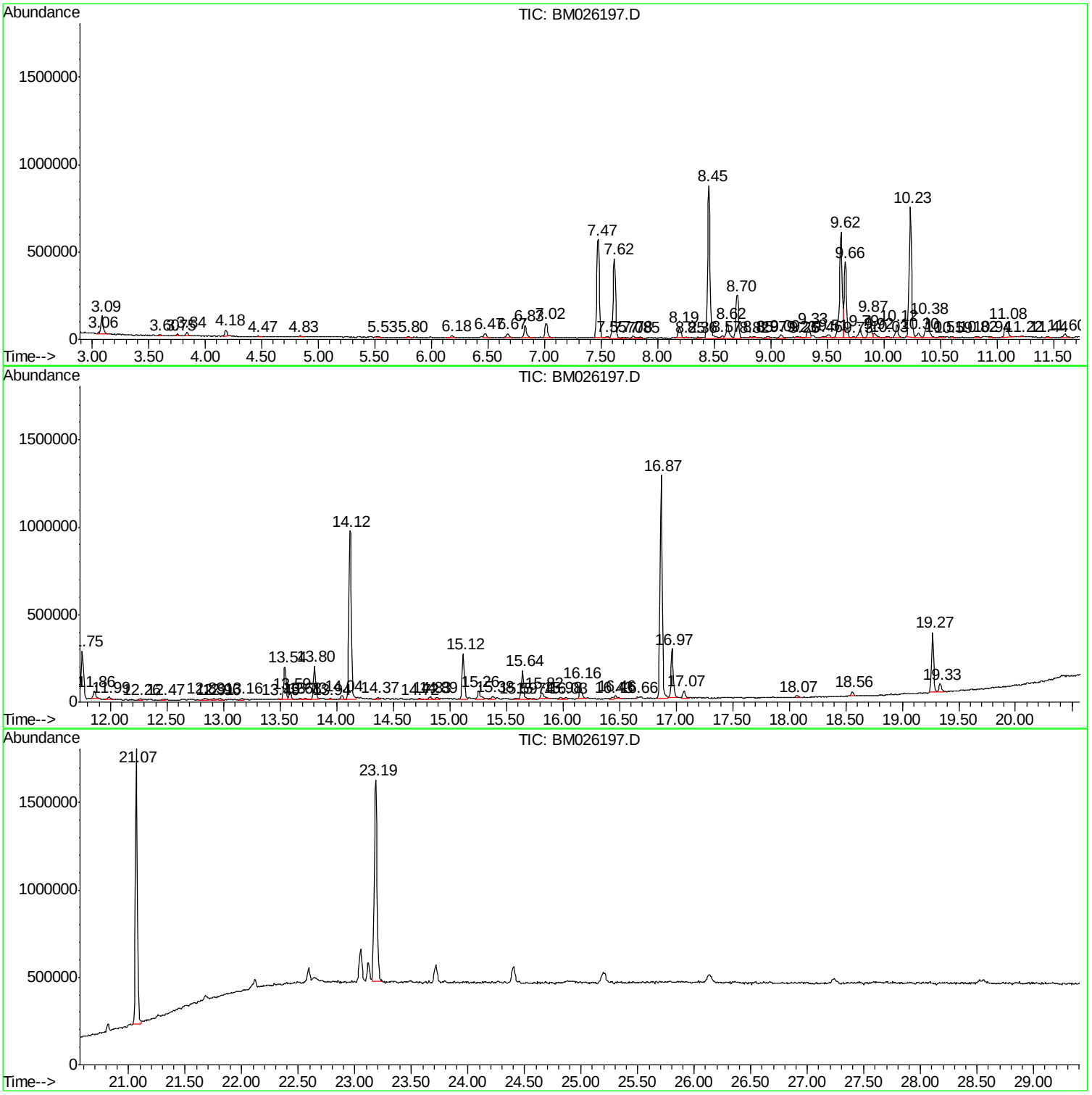
Sum of corrected areas: 19097991

Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM060120\
Data File : BM026197.D
Acq On : 01 Jun 2020 13:16
Operator : JU/CG
Sample : L2820-08DL 5X
Misc :
ALS Vial : 3 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
CB638DL

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\
Data File : BM026197.D
Acq On : 01 Jun 2020 13:16
Operator : JU/CG
Sample : L2820-08DL 5X
Misc :
ALS Vial : 3 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : CB638DL

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

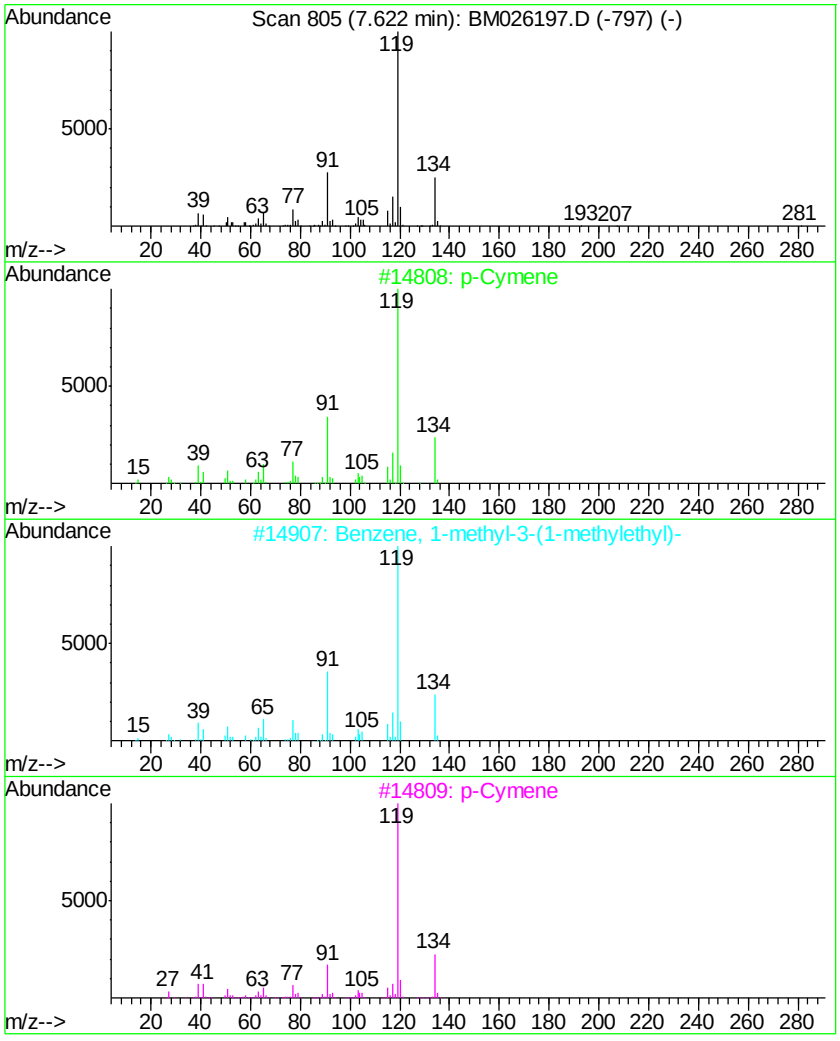
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

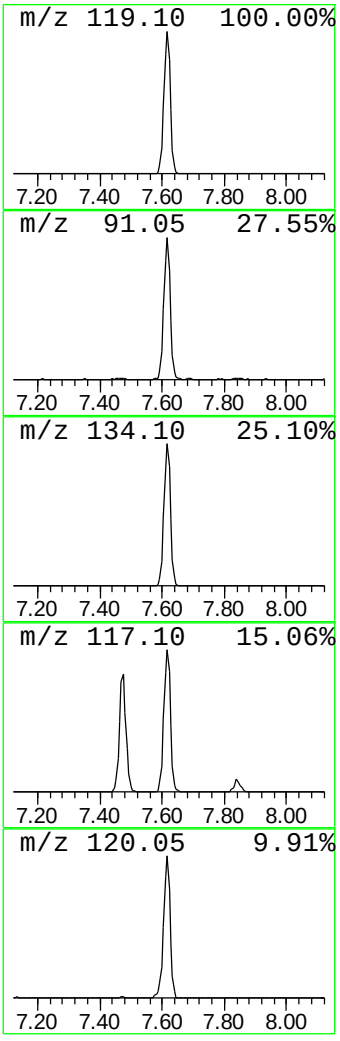
****

Peak Number 1 p-Cymene Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.62	15.13 ng/ul	668810	1,4-Dichlorobenzene-d4	7.47

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	p-Cymene	134	C10H14	000099-87-6	97
2	Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	96
3	p-Cymene	134	C10H14	000099-87-6	94
4	o-Cymene	134	C10H14	000527-84-4	94
5	o-Cymene	134	C10H14	000527-84-4	94





Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM060120\
Data File : BM026197.D
Acq On : 01 Jun 2020 13:16
Operator : JU/CG
Sample : L2820-08DL 5X
Misc :
ALS Vial : 3 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : CB638DL

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

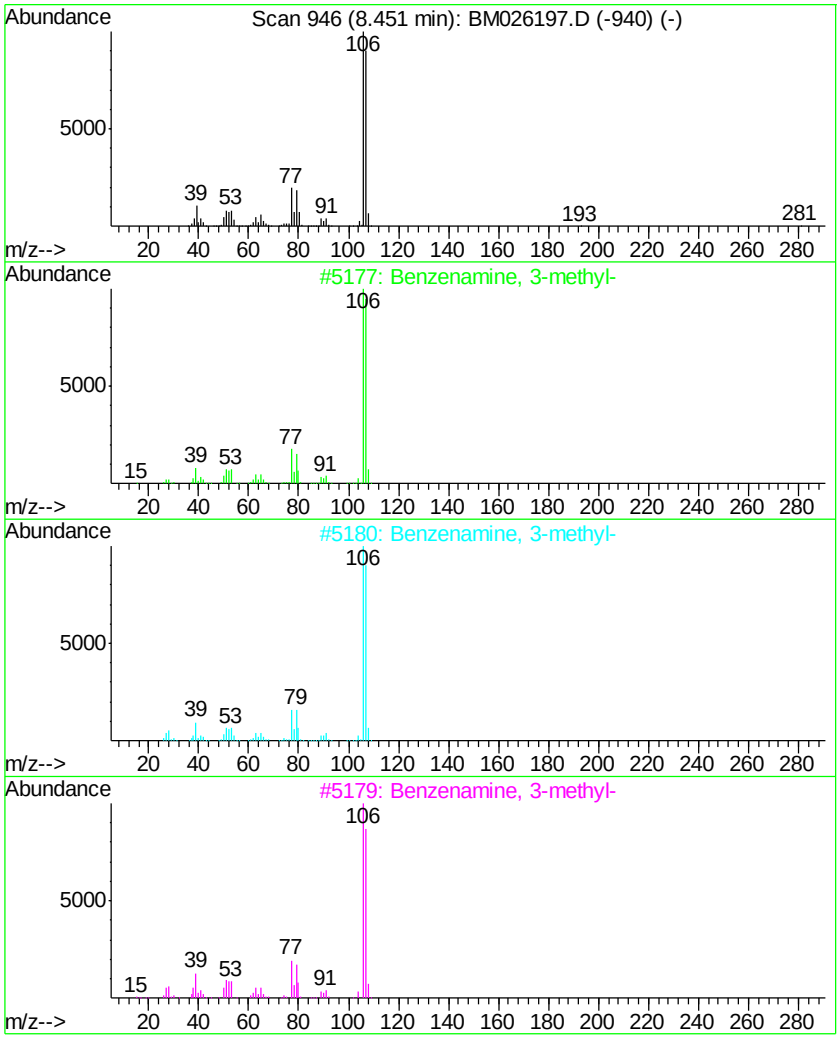
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

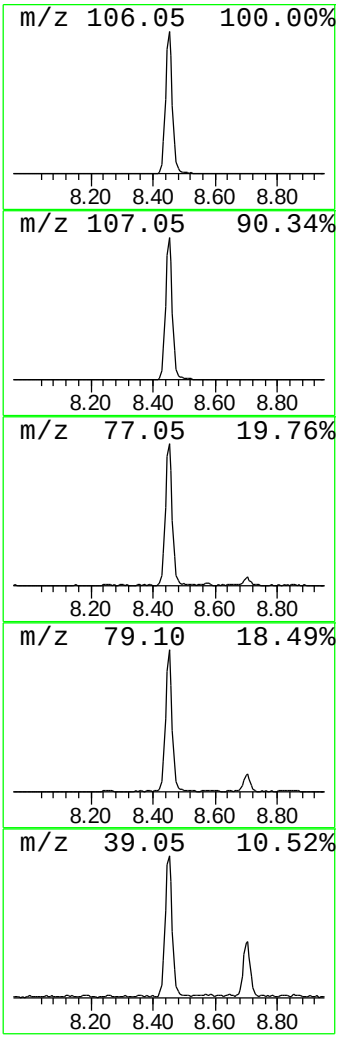
****

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.45	30.70 ng/ul	1356910	1,4-Dichlorobenzene-d4	7.47

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	96
4	Benzenamine, 3-methyl-	107	C7H9N	000108-44-1	95
5	p-Aminotoluene	107	C7H9N	000106-49-0	94





Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM060120\
Data File : BM026197.D
Acq On : 01 Jun 2020 13:16
Operator : JU/CG
Sample : L2820-08DL 5X
Misc :
ALS Vial : 3 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : CB638DL

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

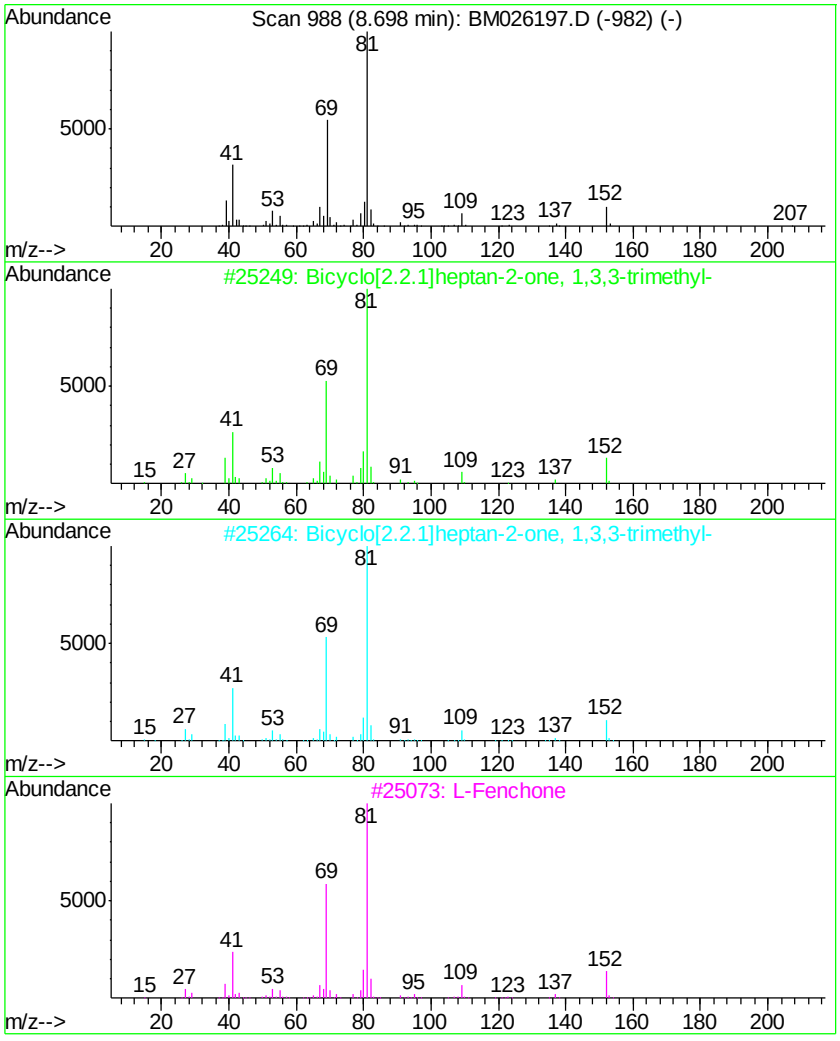
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

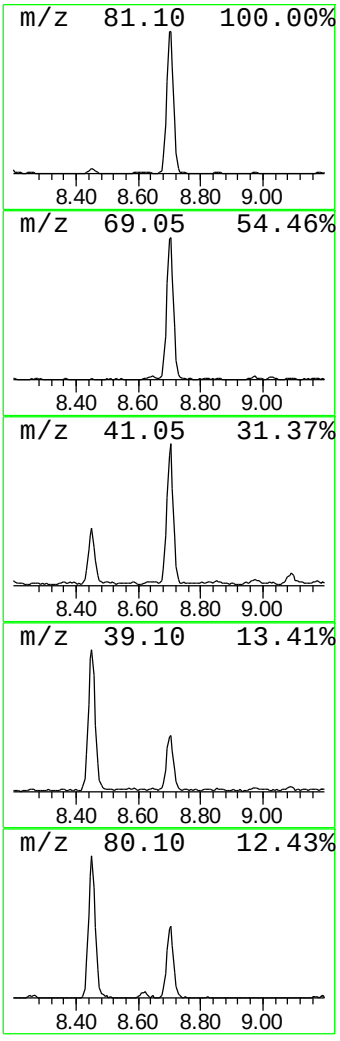
****

Peak Number 3 Bicyclo[2.2.1]heptan-2-one,... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.70	8.86 ng/ul	391840	1,4-Dichlorobenzene-d4	7.47

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	Bicyclo[2.2.1]heptan-2-one, 1,3,...	152	C10H16O	001195-79-5	95
3	L-Fenchone	152	C10H16O	007787-20-4	94
4	L-Fenchone	152	C10H16O	007787-20-4	94
5	Bicyclo[2.2.1]heptan-2-one, 1,3,...	152	C10H16O	001195-79-5	91





Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM060120\
Data File : BM026197.D
Acq On : 01 Jun 2020 13:16
Operator : JU/CG
Sample : L2820-08DL 5X
Misc :
ALS Vial : 3 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : CB638DL

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

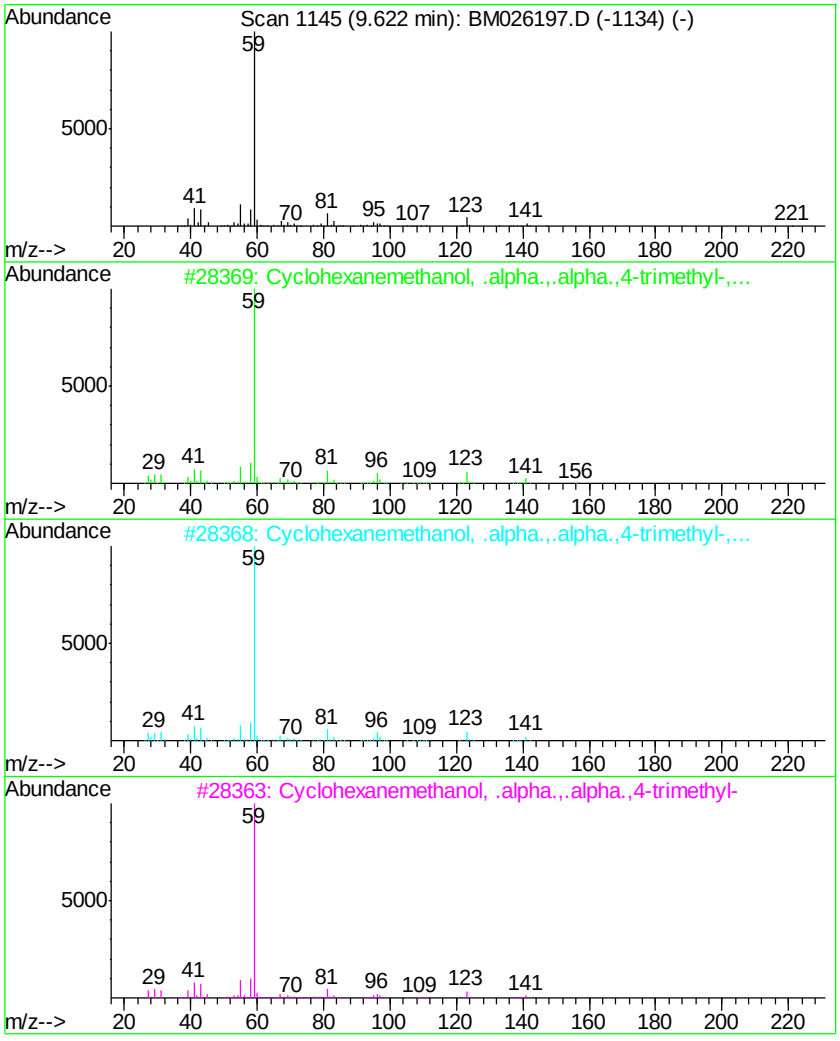
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

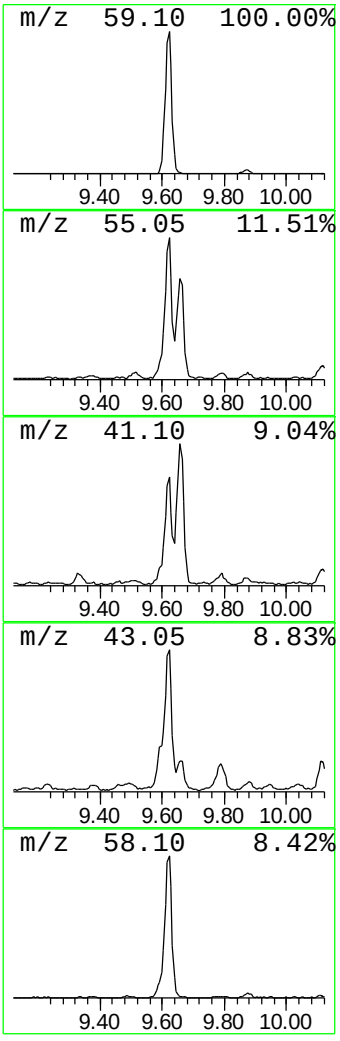
****

Peak Number 4 Cyclohexanemethanol, .alpha... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.62	18.26 ng/ul	1057710	Naphthalene-d8	10.23

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclohexanemethanol, .alpha.,.al...	156	C10H20O	005114-00-1	83
2	Cyclohexanemethanol, .alpha.,.al...	156	C10H20O	007322-63-6	83
3	Cyclohexanemethanol, .alpha.,.al...	156	C10H20O	000498-81-7	74
4	Cyclohexanemethanol, .alpha.,.al...	156	C10H20O	000498-81-7	47
5	o-Menthan-8-ol	156	C10H20O	343855-44-7	40





Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM060120\
Data File : BM026197.D
Acq On : 01 Jun 2020 13:16
Operator : JU/CG
Sample : L2820-08DL 5X
Misc :
ALS Vial : 3 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : CB638DL

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

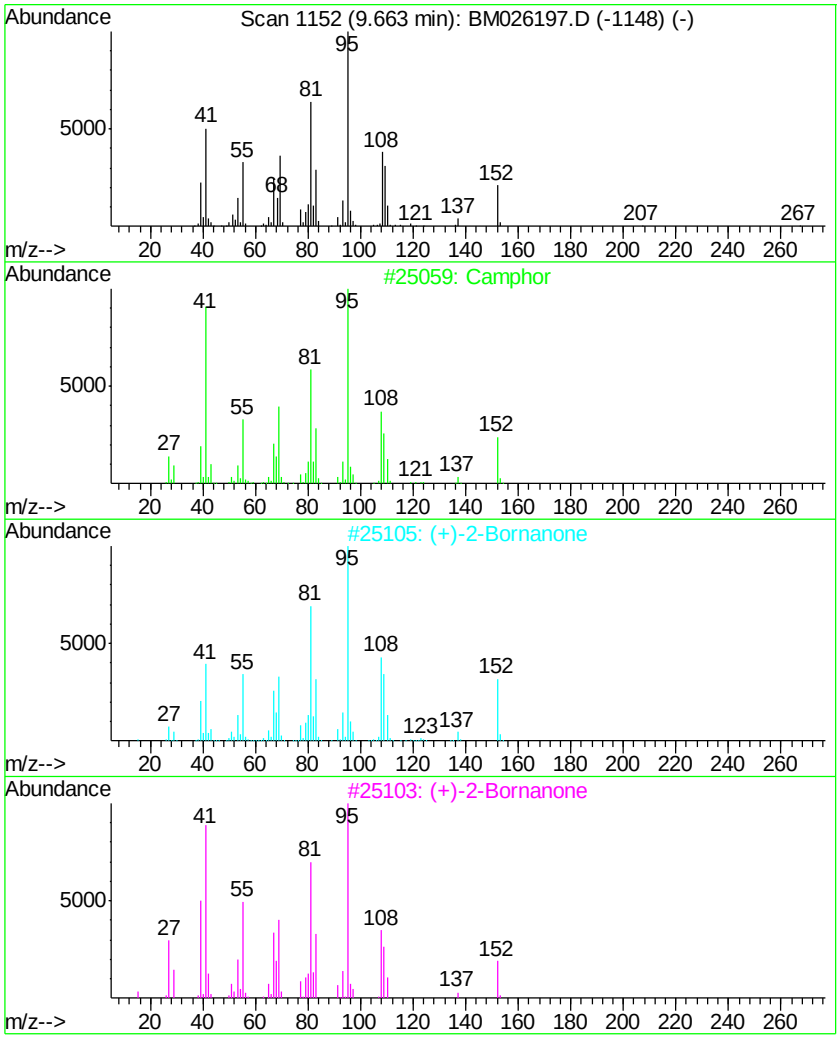
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

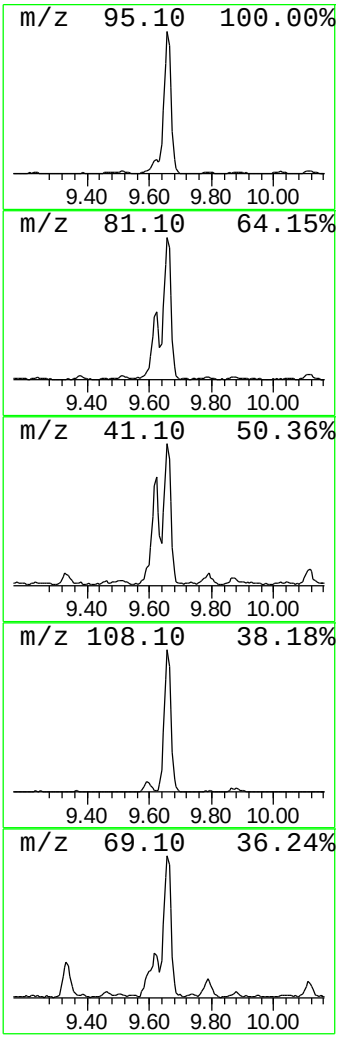
****

Peak Number 5 Camphor Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.66	11.20 ng/ul	648572	Naphthalene-d8	10.23

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Camphor	152	C10H16O	000076-22-2	95
2	(+)-2-Bornanone	152	C10H16O	000464-49-3	94
3	(+)-2-Bornanone	152	C10H16O	000464-49-3	94
4	Camphor	152	C10H16O	000076-22-2	93
5	Bicyclo[2.2.1]heptan-2-one, 1,7,...	152	C10H16O	000464-48-2	92





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Acq On : 01 Jun 2020 13:16
Operator : JU/CG
Sample : L2820-08DL 5X
Misc :
ALS Vial : 3 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : CB638DL

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

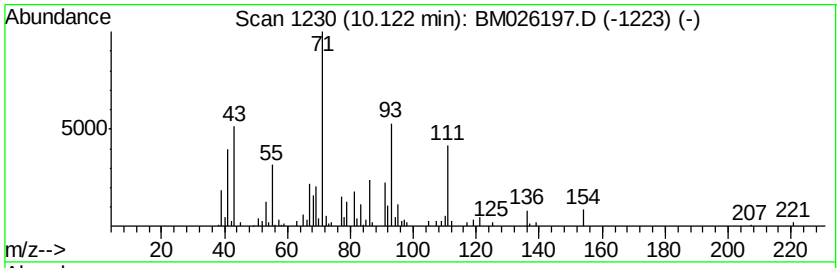
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

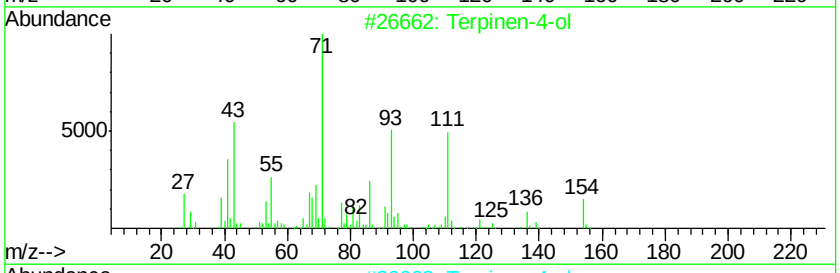
****

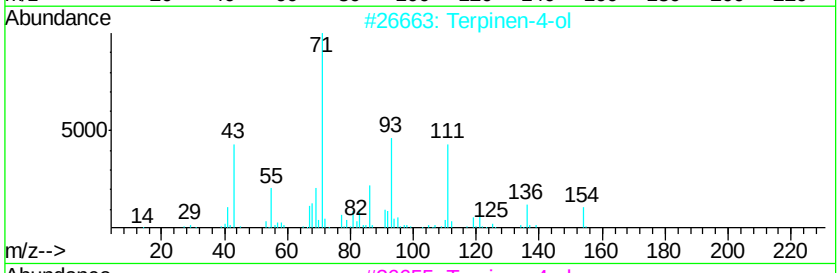
Peak Number 6 Terpinen-4-ol Concentration Rank 9

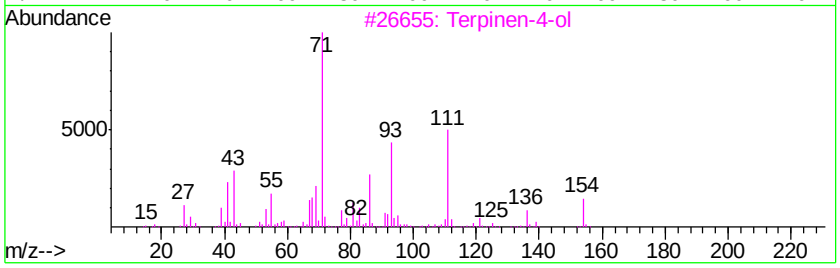
R.T.	EstConc	Area	Relative to ISTD	R.T.
10.12	2.06 ng/ul	119221	Naphthalene-d8	10.23

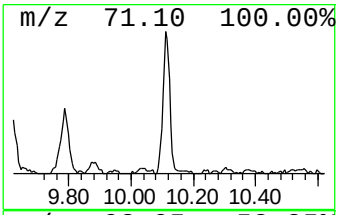
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Terpinen-4-ol	154	C10H18O	000562-74-3	95
2	Terpinen-4-ol	154	C10H18O	000562-74-3	91
3	Terpinen-4-ol	154	C10H18O	000562-74-3	81
4	3-Cyclohexen-1-ol, 4-methyl-1-(1...	154	C10H18O	020126-76-5	81
5	Terpinen-4-ol	154	C10H18O	000562-74-3	68

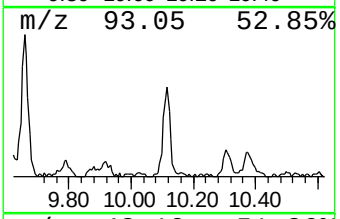


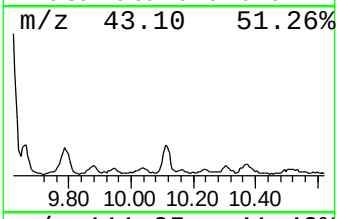


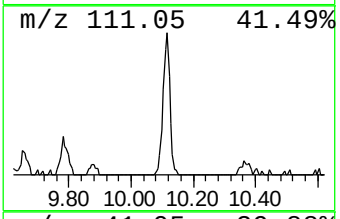


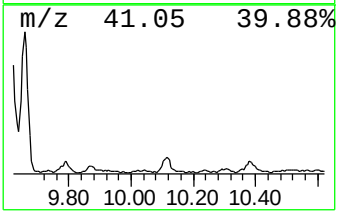












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Data File : BM026197.D
Acq On : 01 Jun 2020 13:16
Operator : JU/CG
Sample : L2820-08DL 5X
Misc :
ALS Vial : 3 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : CB638DL

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

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

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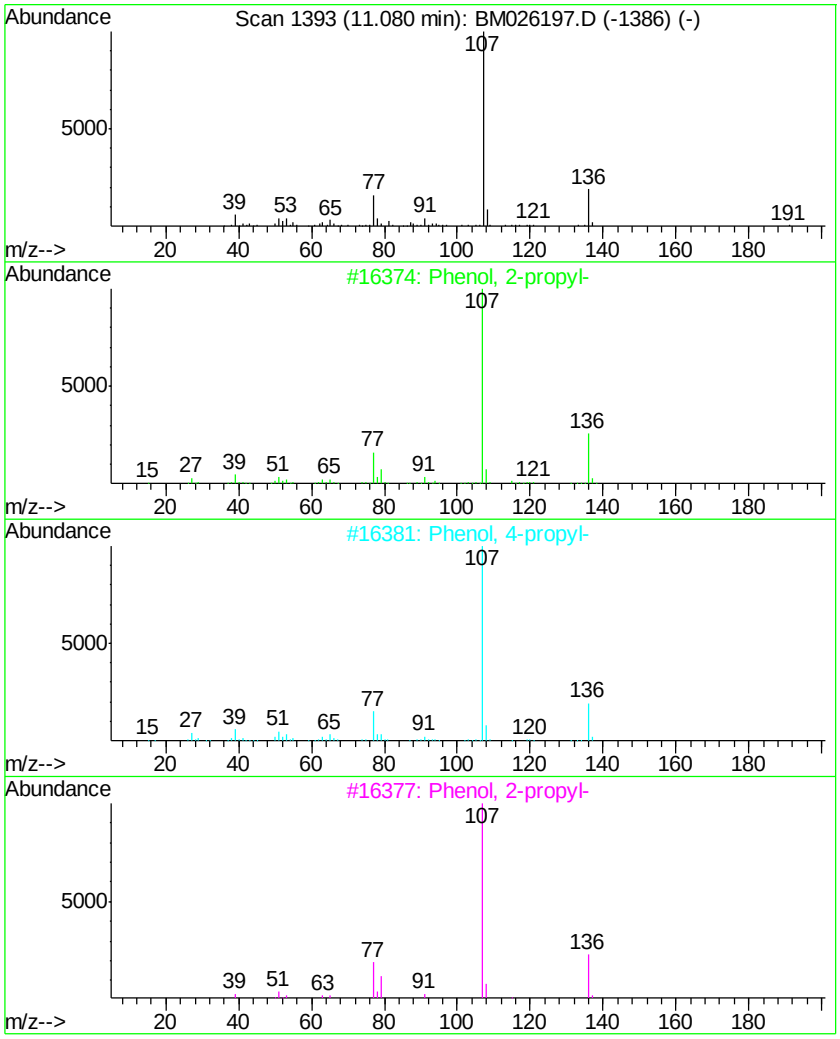
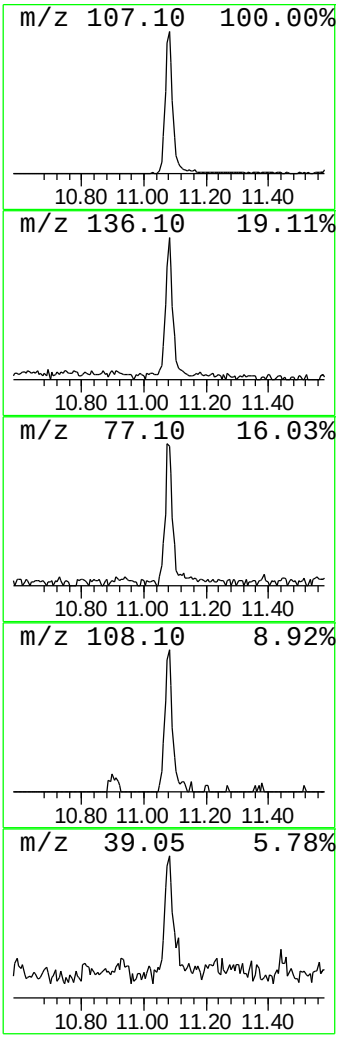
Peak Number 7 Phenol, 2-propyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.08	2.57 ng/ul	149150	Naphthalene-d8	10.23

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, 2-propyl-	136	C9H12O	000644-35-9	91
2	Phenol, 4-propyl-	136	C9H12O	000645-56-7	90
3	Phenol, 2-propyl-	136	C9H12O	000644-35-9	86
4	Phenol, 4-propyl-	136	C9H12O	000645-56-7	80
5	Formic acid, 2-propylphenyl ester	164	C10H12O2	1000368-96-6	78





Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM060120\
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Operator : JU/CG
Sample : L2820-08DL 5X
Misc :
ALS Vial : 3 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : CB638DL

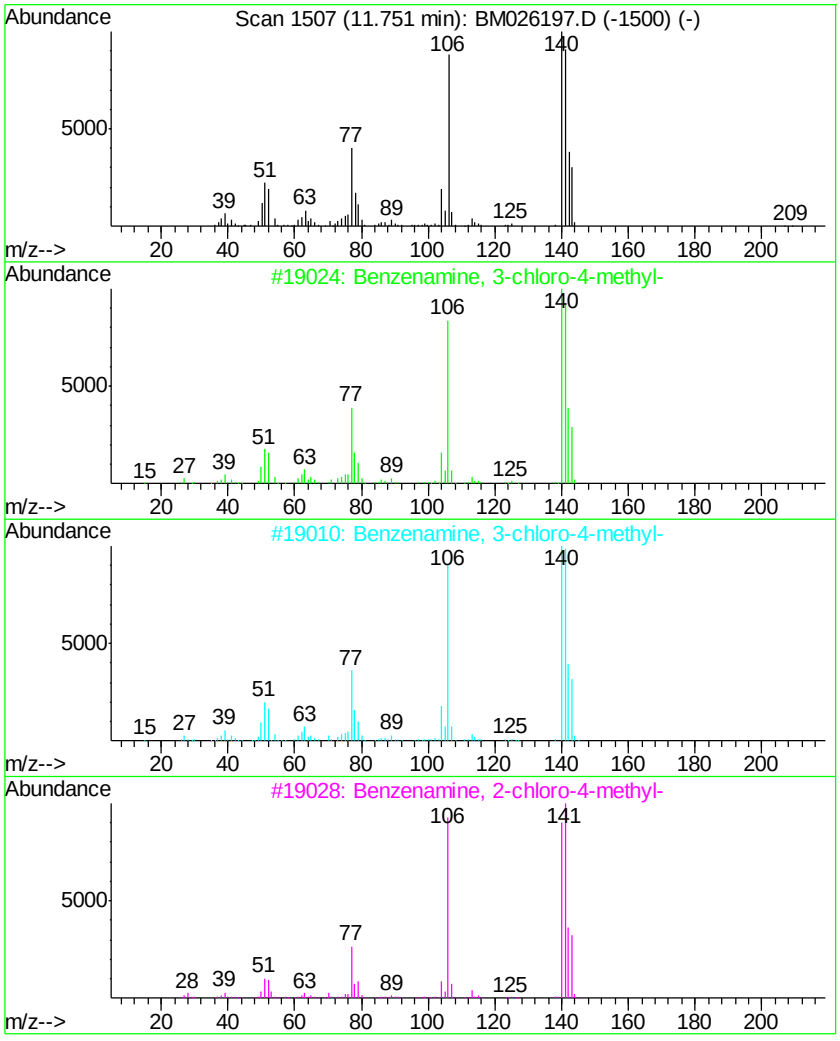
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM051320MA.M
Quant Title : SVOA CALIBRATION

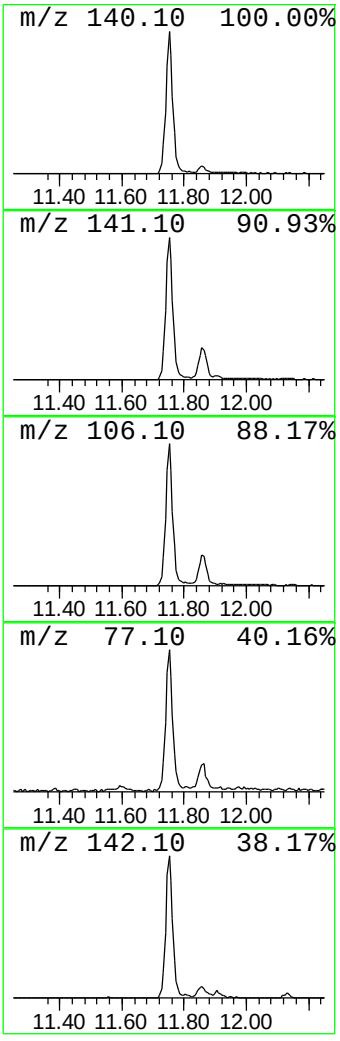
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 8 Benzenamine, 3-chloro-4-met... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.75	7.71 ng/ul	446639	Naphthalene-d8	10.23

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzenamine, 3-chloro-4-methyl-	141	C7H8ClN	000095-74-9	98
2	Benzenamine, 3-chloro-4-methyl-	141	C7H8ClN	000095-74-9	98
3	Benzenamine, 2-chloro-4-methyl-	141	C7H8ClN	000615-65-6	95
4	Benzenamine, 2-chloro-4-methyl-	141	C7H8ClN	000615-65-6	95
5	Benzenamine, 3-chloro-4-methyl-	141	C7H8ClN	000095-74-9	93





Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM060120\
Data File : BM026197.D
Acq On : 01 Jun 2020 13:16
Operator : JU/CG
Sample : L2820-08DL 5X
Misc :
ALS Vial : 3 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : CB638DL

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

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

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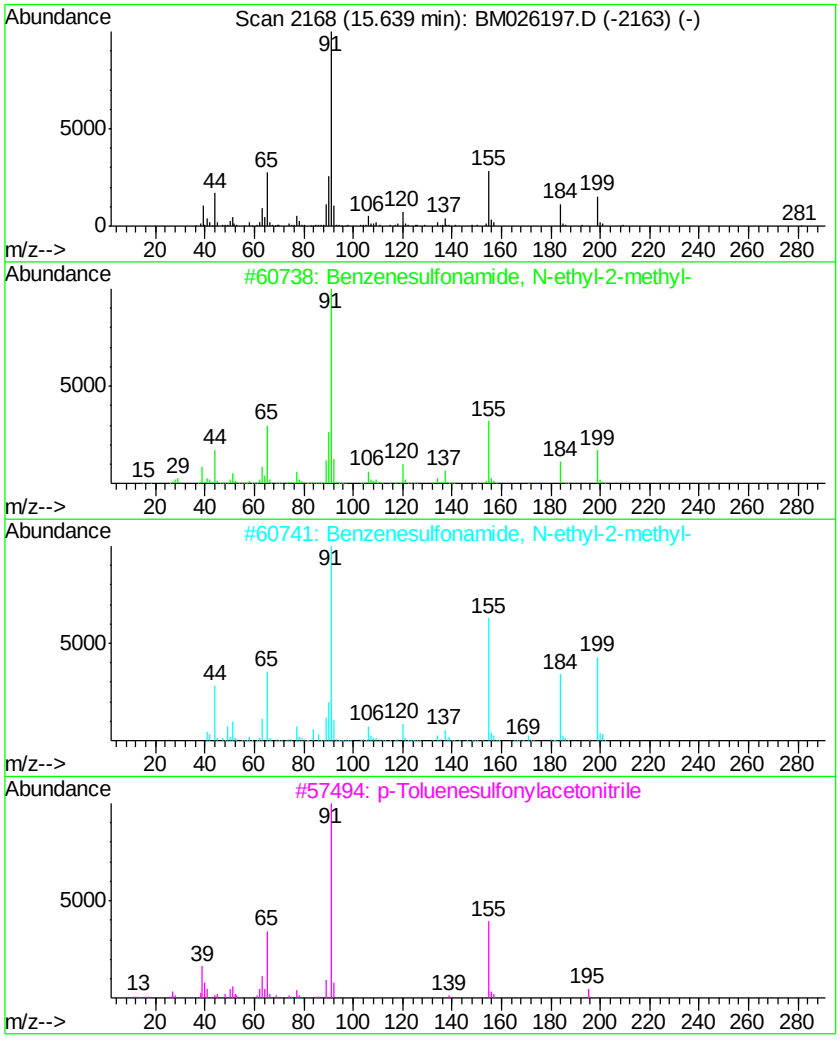
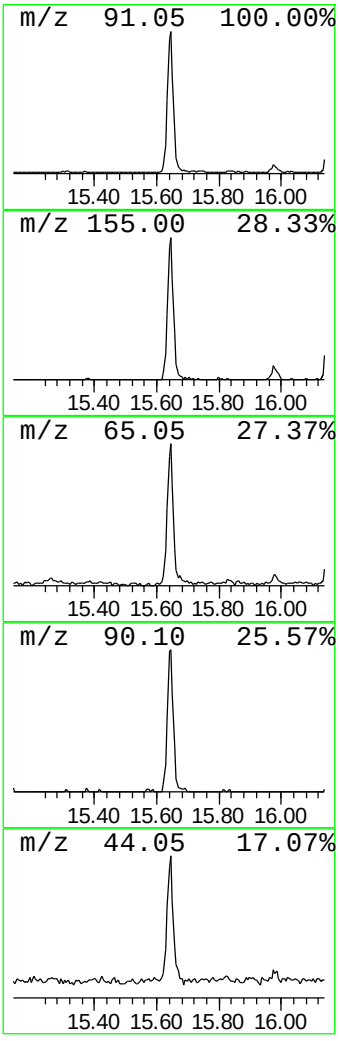
Peak Number 9 Benzenesulfonamide, N-ethyl... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.64	2.80 ng/ul	243835	Phenanthrene-d10	16.87

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzenesulfonamide, N-ethyl-2-me...	199	C9H13NO2S	001077-56-1	98
2	Benzenesulfonamide, N-ethyl-2-me...	199	C9H13NO2S	001077-56-1	64
3	p-Toluenesulfonylacetonitrile	195	C9H9NO2S	005697-44-9	50
4	Ethyl(E/Z)-.alpha.-cyano-4-nitro...	416	C19H16N2O7S	1000287-26-8	50
5	4-Toluenesulfonylmethyl isocyanide	195	C9H9NO2S	036635-61-7	47





Tentatively Identified Compound (LSC) summary

Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM060120\
Data File : BM026197.D
Acq On : 01 Jun 2020 13:16
Operator : JU/CG
Sample : L2820-08DL 5X
Misc :
ALS Vial : 3 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : CB638DL

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

					Internal Standard			
TIC Top Hit name	RT	EstConc	Units	Response	#	RT	Resp	Conc
p-Cymene	7.62	15.1	ng/ul	668810	1	7.47	884035	20.0
Benzenamine, 3-me...	8.45	30.7	ng/ul	1356910	1	7.47	884035	20.0
Bicyclo[2.2.1]hep...	8.70	8.9	ng/ul	391840	1	7.47	884035	20.0
Cyclohexanemethan...	9.62	18.3	ng/ul	1057710	2	10.23	1158510	20.0
Camphor	9.66	11.2	ng/ul	648572	2	10.23	1158510	20.0
Terpinen-4-ol	10.12	2.1	ng/ul	119221	2	10.23	1158510	20.0
Phenol, 2-propyl-	11.08	2.6	ng/ul	149150	2	10.23	1158510	20.0
Benzenamine, 3-ch...	11.75	7.7	ng/ul	446639	2	10.23	1158510	20.0
Benzenesulfonamid...	15.64	2.8	ng/ul	243835	4	16.87	1743720	20.0