LSC Area Percent Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM052920\
Data File : BM026192.D
Acq On : 30 May 2020 03:47
Operator : MA/SJ
Sample : L2820-13
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : CB640

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	3.063	27	30	33	rBV	45774	69133	0.80%	0.093%
2	3.099	33	36	48	rVB	80033	142969	1.65%	0.193%
3	4.187	216	221	234	rVB	112240	190131	2.19%	0.257%
4	4.428	257	262	267	rBV	38619	64161	0.74%	0.087%
5	4.846	326	333	338	rBV	107029	158854	1.83%	0.215%
6	5.928	511	517	524	rVB	38509	61632	0.71%	0.083%
7	5.993	524	528	533	rBV	42782	69883	0.80%	0.094%
8	6.187	556	561	566	rVB	67587	102232	1.18%	0.138%
9	6.481	606	611	615	rVV2	45759	71918	0.83%	0.097%
10	6.675	639	644	655	rVB2	201164	366771	4.22%	0.496%
11	6.828	660	670	677	rBV	529804	814824	9.38%	1.102%
12	7.016	696	702	712	rVV	610697	966066	11.12%	1.306%
13	7.363	755	761	769	rVV	21675	59677	0.69%	0.081%
14	7.475	769	780	789	rVV	673022	1159372	13.34%	1.567%
15	7.551	789	793	797	rVV2	54780	92160	1.06%	0.125%
16	7.616	799	804	815	rVB	312826	482586	5.55%	0.652%
17	7.845	836	843	849	rVB	116619	189771	2.18%	0.257%
18	8.192	897	902	907	rVV	483981	810642	9.33%	1.096%
19	8.257	907	913	927	rVV	5655181	8689613	100.00%	11.748%
20	8.457	941	947	953	rVV	263484	409117	4.71%	0.553%
21	8.575	961	967	970	rBV	72552	119170	1.37%	0.161%
22	8.622	970	975	985	rVV	665343	1108740	12.76%	1.499%
23	8.710	985	990	997	rVB	152703	234725	2.70%	0.317%
24	8.975	1028	1035	1042	rBV5	56766	109640	1.26%	0.148%
25	9.098	1049	1056	1062	rBV3	80090	142090	1.64%	0.192%
26	9.334	1090	1096	1109	rBV	551931	927621	10.68%	1.254%
27	9.622	1138	1145	1147	rVV2	109990	190000	2.19%	0.257%
28	9.657	1147	1151	1160	rVB2	337247	576580	6.64%	0.779%
29	9.792	1161	1174	1180	rVB3	124107	308397	3.55%	0.417%
30	9.869	1180	1187	1198	rBV	884426	1518793	17.48%	2.053%
31	10.022	1208	1213	1222	rBV6	32739	88191	1.01%	0.119%
32	10.116	1222	1229	1236	rVV	177342	307922	3.54%	0.416%
33	10.233	1240	1249	1254	rVV	902038	1508848	17.36%	2.040%
34	10.286	1254	1258	1266	rVB	210586	395556	4.55%	0.535%

LSC Area Percent Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM052920\
Data File : BM026192.D
Acq On : 30 May 2020 03:47
Operator : MA/SJ
Sample : L2820-13
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : CB640

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	10.380	1266	1274	1286	rBV	762991	1357915	15.63%	1.836%
36	10.598	1307	1311	1317	rVB5	33828	56641	0.65%	0.077%
37	10.816	1339	1348	1356	rBV4	47462	127674	1.47%	0.173%
38	11.080	1382	1393	1405	rBV5	73365	209835	2.41%	0.284%
39	11.445	1450	1455	1465	rBV8	28358	71253	0.82%	0.096%
40	11.604	1473	1482	1488	rBV	212857	371324	4.27%	0.502%
41	11.751	1501	1507	1512	rBV4	96430	171042	1.97%	0.231%
42	11.822	1512	1519	1523	rBV7	27263	71519	0.82%	0.097%
43	11.904	1528	1533	1539	rVV	236354	423702	4.88%	0.573%
44	11.969	1539	1544	1552	rVV4	59762	141377	1.63%	0.191%
45	12.092	1561	1565	1567	rVV2	93471	134381	1.55%	0.182%
46	12.127	1567	1571	1578	rVB	204284	348461	4.01%	0.471%
47	12.463	1625	1628	1634	rVB7	32105	61979	0.71%	0.084%
48	12.969	1703	1714	1722	rVB	560398	1010692	11.63%	1.366%
49	13.169	1741	1748	1755	rVB7	79352	167622	1.93%	0.227%
50	13.274	1760	1766	1769	rBV6	71285	144651	1.66%	0.196%
51	13.363	1778	1781	1788	rVB7	36871	65741	0.76%	0.089%
52	13.439	1789	1794	1799	rBV	83887	157791	1.82%	0.213%
53	13.492	1799	1803	1807	rVV3	70250	103133	1.19%	0.139%
54	13.545	1807	1812	1816	rVV	1600811	2071642	23.84%	2.801%
55	13.592	1816	1820	1824	rVB	234012	292300	3.36%	0.395%
56	13.680	1831	1835	1837	rBV3	47969	67990	0.78%	0.092%
57	13.810	1850	1857	1866	rBV	1587838	2476998	28.51%	3.349%
58	13.951	1875	1881	1889	rBV2	490079	860011	9.90%	1.163%
59	14.051	1893	1898	1903	rVV2	153226	238734	2.75%	0.323%
60	14.121	1903	1910	1915	rVV2	2020647	3087522	35.53%	4.174%
61	14.186	1915	1921	1927	rVV	1200333	1813794	20.87%	2.452%
62	14.251	1927	1932	1939	rVB	362895	506026	5.82%	0.684%
63	14.363	1947	1951	1956	rBV	138776	185512	2.13%	0.251%
64	14.521	1973	1978	1988	rVB	460433	675687	7.78%	0.913%
65	14.680	2001	2005	2010	rVB	167921	210327	2.42%	0.284%
66	14.886	2032	2040	2050	rBV2	1195485	2293640	26.40%	3.101%
67	15.121	2074	2080	2085	rBV	2133354	2933836	33.76%	3.966%
68	15.174	2085	2089	2094	rVB	832172	1150423	13.24%	1.555%
69	15.251	2098	2102	2107	rBV	569331	805511	9.27%	1.089%
70	15.298	2108	2110	2114	rVB3	42354	54634	0.63%	0.074%
71	15.386	2121	2125	2130	rBV2	445528	566686	6.52%	0.766%

LSC Area Percent Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM052920\
Data File : BM026192.D
Acq On : 30 May 2020 03:47
Operator : MA/SJ
Sample : L2820-13
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : CB640

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	15.474	2136	2140	2147	rBV2	76454	137866	1.59%	0.186%
73	15.651	2162	2170	2178	rVB	1310759	1821448	20.96%	2.462%
74	15.821	2194	2199	2211	rBV	530727	919380	10.58%	1.243%
75	15.980	2222	2226	2230	rBV	74783	111262	1.28%	0.150%
76	16.162	2252	2257	2264	rVV	734883	1005683	11.57%	1.360%
77	16.227	2264	2268	2271	rBV2	95502	120693	1.39%	0.163%
78	16.439	2301	2304	2311	rVB	113565	168847	1.94%	0.228%
79	16.692	2343	2347	2353	rVB3	178678	256843	2.96%	0.347%
80	16.868	2372	2377	2381	rVV	1545750	2074000	23.87%	2.804%
81	16.909	2381	2384	2389	rVB	1221434	1612498	18.56%	2.180%
82	16.968	2389	2394	2399	rBV	2413392	3236027	37.24%	4.375%
83	17.068	2408	2411	2417	rVB2	128699	180282	2.07%	0.244%
84	17.280	2442	2447	2452	rVB	531690	704993	8.11%	0.953%
85	17.386	2461	2465	2469	rBV4	50475	77425	0.89%	0.105%
86	17.433	2470	2473	2478	rBV2	46830	80363	0.92%	0.109%
87	17.715	2518	2521	2524	rBV	55702	65281	0.75%	0.088%
88	17.798	2531	2535	2543	rBV3	87834	140948	1.62%	0.191%
89	17.874	2545	2548	2552	rVB2	71349	92074	1.06%	0.124%
90	17.939	2553	2559	2564	rBV2	164119	268044	3.08%	0.362%
91	18.562	2662	2665	2672	rVB4	90786	138774	1.60%	0.188%
92	18.933	2725	2728	2733	rVB	318605	392756	4.52%	0.531%
93	19.027	2741	2744	2748	rVB	154464	171184	1.97%	0.231%
94	19.268	2780	2785	2793	rBV	2616956	3796355	43.69%	5.132%
95	19.339	2793	2797	2807	rVB	426649	630482	7.26%	0.852%
96	19.756	2866	2868	2874	rVB	99713	116699	1.34%	0.158%
97	20.815	3044	3048	3054	rBV	597324	711784	8.19%	0.962%
98	21.068	3087	3091	3099	rVB	1918509	2320768	26.71%	3.137%
99	23.050	3423	3428	3435	rVB	1405974	2293986	26.40%	3.101%
100	23.186	3445	3451	3459	rVB	1358880	2326689	26.78%	3.145%

Sum of corrected areas: 73969225

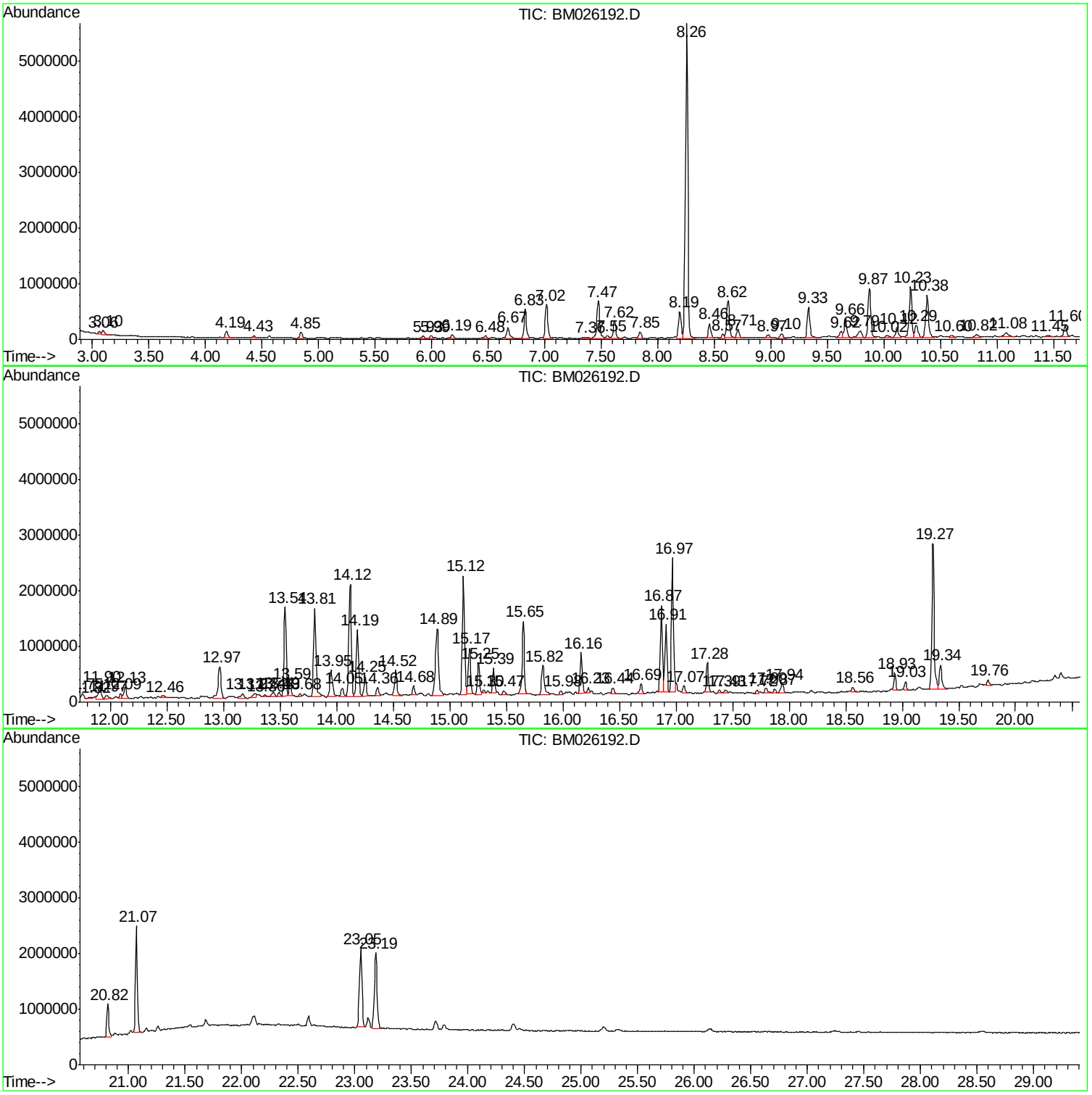
LSC Report - Integrated Chromatogram

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM052920\
Data File : BM026192.D
Acq On : 30 May 2020 03:47
Operator : MA/SJ
Sample : L2820-13
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : CB640

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



Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM052920\
Data File : BM026192.D
Acq On : 30 May 2020 03:47
Operator : MA/SJ
Sample : L2820-13
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : CB640

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

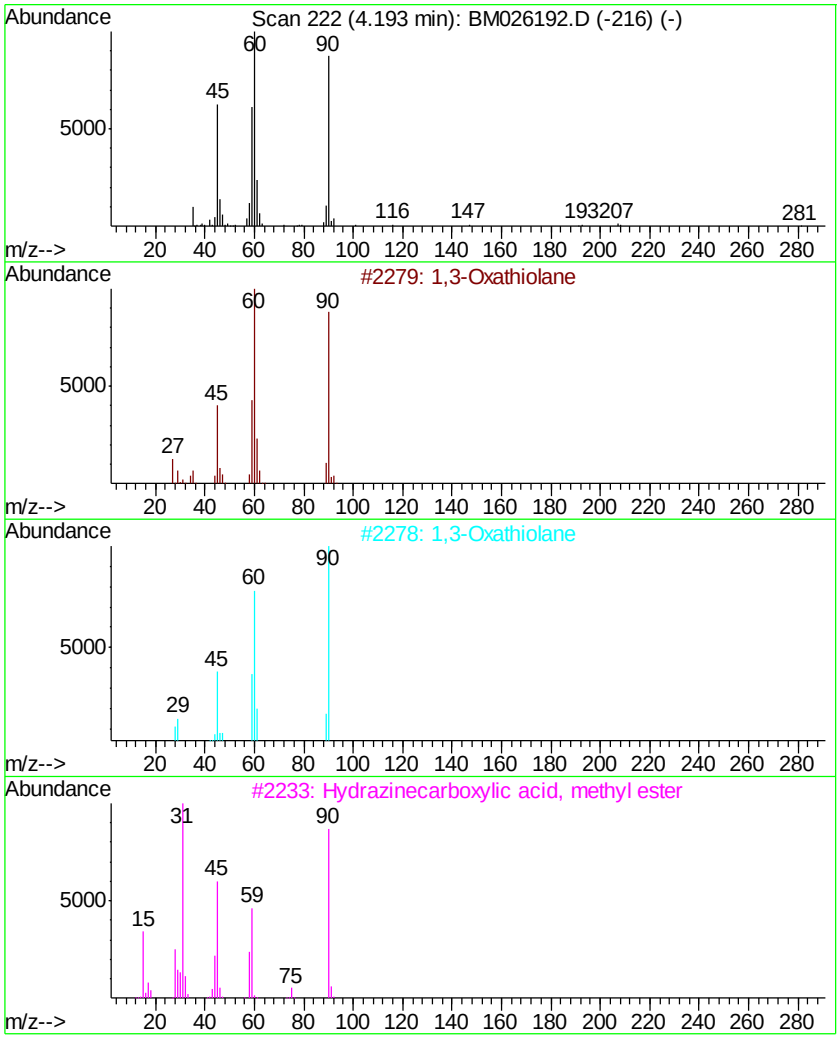
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

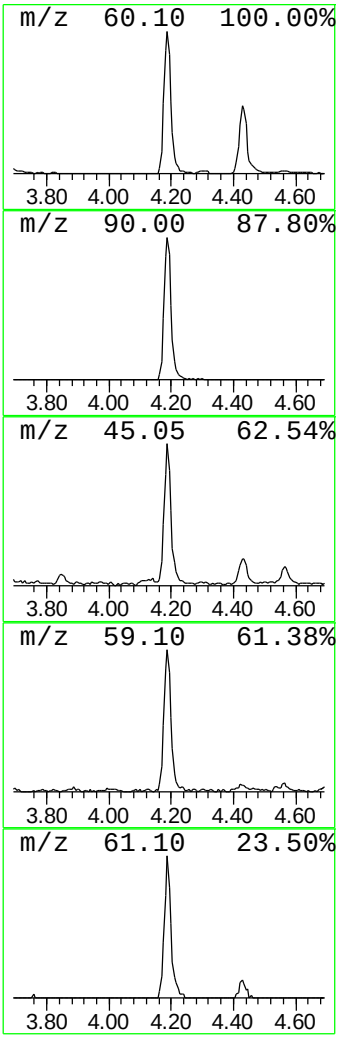
****

Peak Number 1 1,3-Oxathiolane Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.19	3.28 ng/ul	190131	1,4-Dichlorobenzene-d4	7.47

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,3-Oxathiolane	90	C3H6OS	002094-97-5	90
2	1,3-Oxathiolane	90	C3H6OS	002094-97-5	74
3	Hydrazinecarboxylic acid, methyl...	90	C2H6N2O2	006294-89-9	38
4	Hydrazinecarboxylic acid, methyl...	90	C2H6N2O2	006294-89-9	38
5	Heptanoic acid, 2-hydroxy-, meth...	160	C8H16O3	054340-91-9	36





Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM052920\
Data File : BM026192.D
Acq On : 30 May 2020 03:47
Operator : MA/SJ
Sample : L2820-13
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : CB640

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

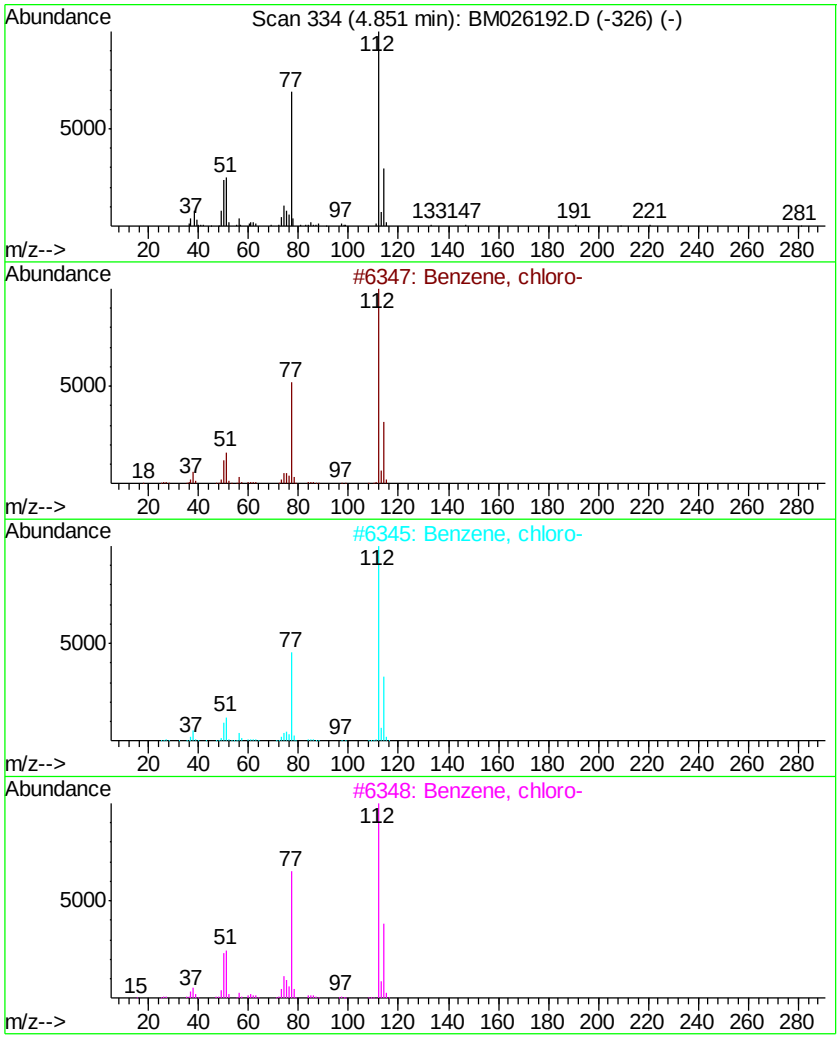
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

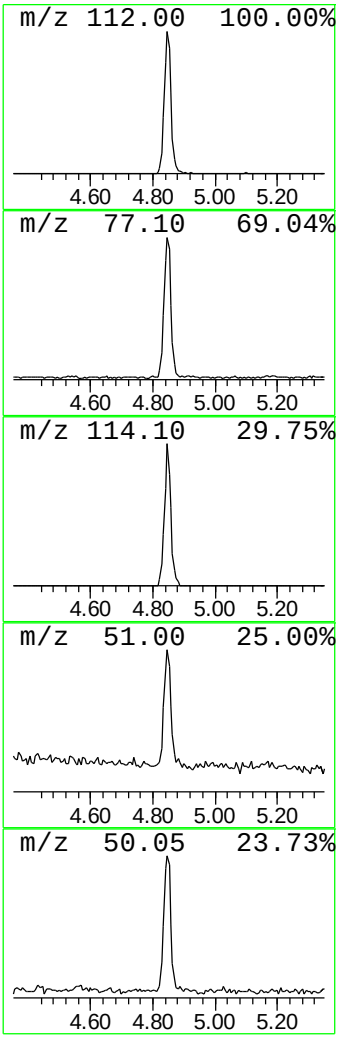
****

Peak Number 2 Benzene, chloro- Concentration Rank 21

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.85	2.74 ng/ul	158854	1,4-Dichlorobenzene-d4	7.47

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, chloro-	112	C6H5Cl	000108-90-7	95
2	Benzene, chloro-	112	C6H5Cl	000108-90-7	95
3	Benzene, chloro-	112	C6H5Cl	000108-90-7	94
4	Benzene, chloro-	112	C6H5Cl	000108-90-7	87
5	2-Chloro-3-nitropyridine	158	C5H3ClN2O2	005470-18-8	28





Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM052920\
Data File : BM026192.D
Acq On : 30 May 2020 03:47
Operator : MA/SJ
Sample : L2820-13
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : CB640

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

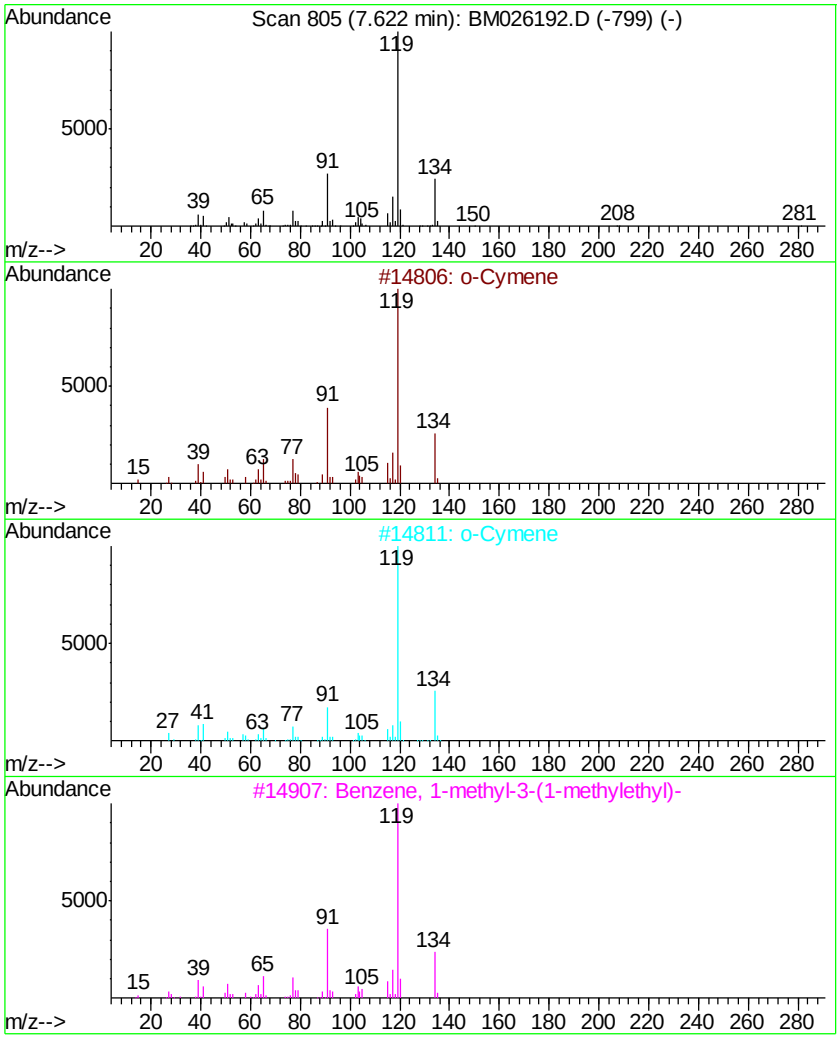
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

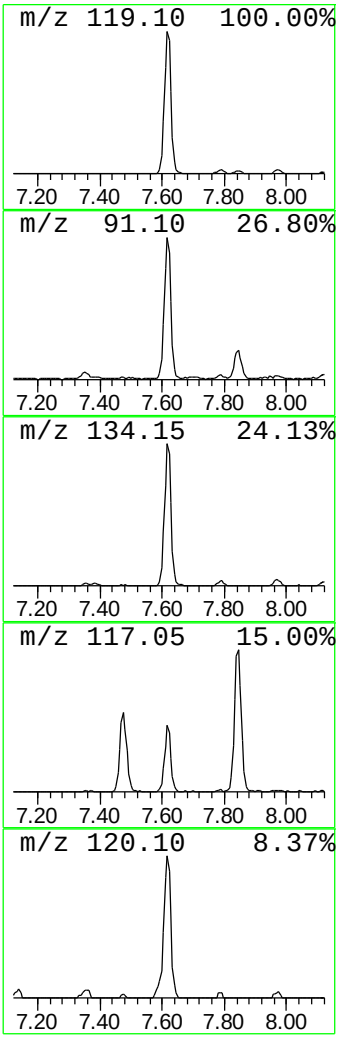
****

Peak Number 3 o-Cymene Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.62	8.32 ng/ul	482586	1,4-Dichlorobenzene-d4	7.47

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	o-Cymene	134	C10H14	000527-84-4	96
2	o-Cymene	134	C10H14	000527-84-4	94
3	Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	94
4	o-Cymene	134	C10H14	000527-84-4	94
5	o-Cymene	134	C10H14	000527-84-4	91





Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM052920\
Data File : BM026192.D
Acq On : 30 May 2020 03:47
Operator : MA/SJ
Sample : L2820-13
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : CB640

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

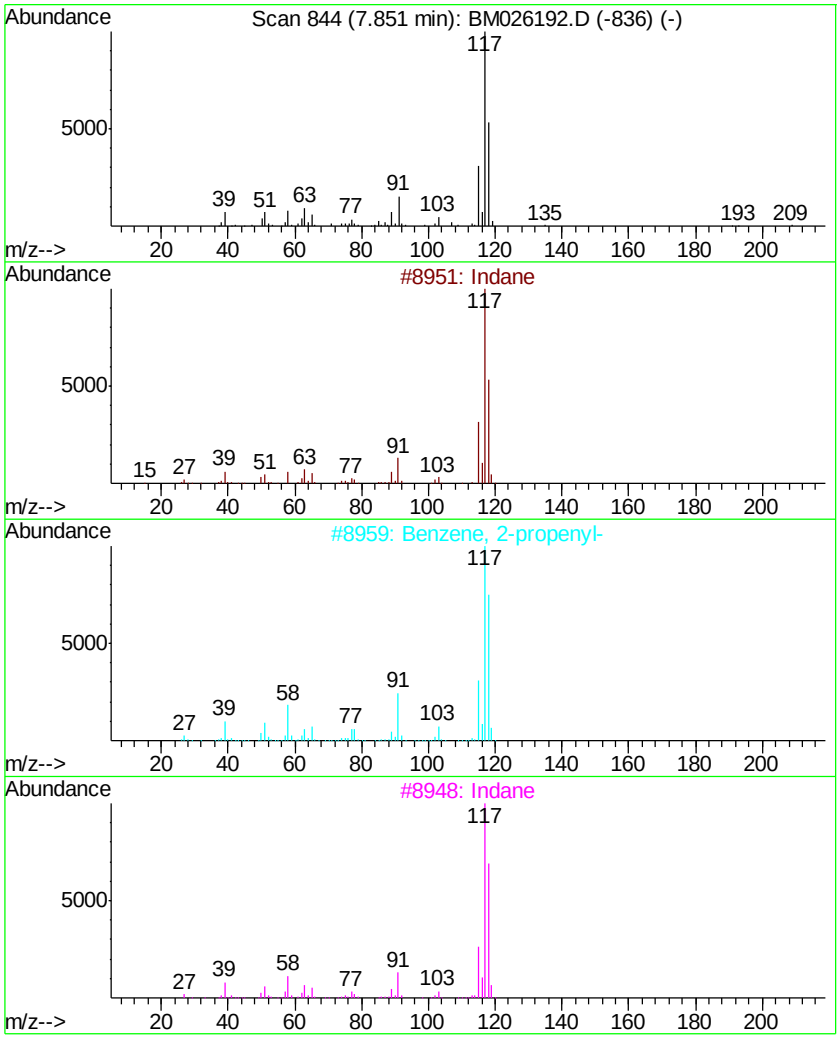
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

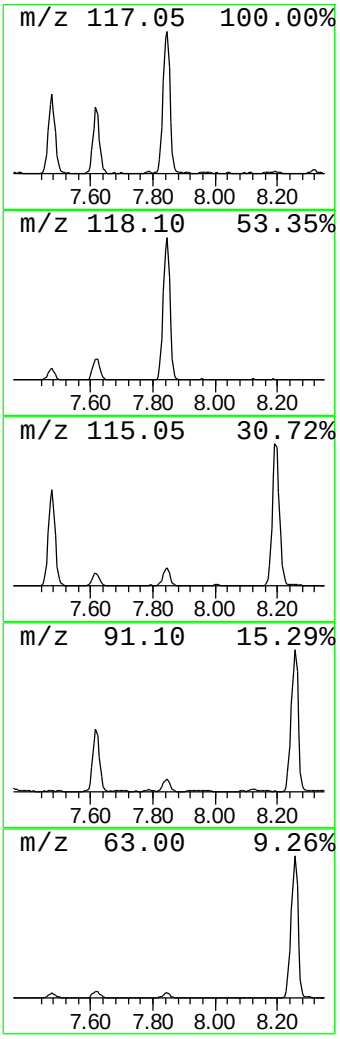
****

Peak Number 4 Indane Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.85	3.27 ng/ul	189771	1,4-Dichlorobenzene-d4	7.47

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indane	118	C9H10	000496-11-7	81
2	Benzene, 2-propenyl-	118	C9H10	000300-57-2	64
3	Indane	118	C9H10	000496-11-7	64
4	Indane	118	C9H10	000496-11-7	60
5	1H-Indene-5-carboxylic acid, 2,3...	162	C10H10O2	1000337-61-3	59





Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM052920\
Data File : BM026192.D
Acq On : 30 May 2020 03:47
Operator : MA/SJ
Sample : L2820-13
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : CB640

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

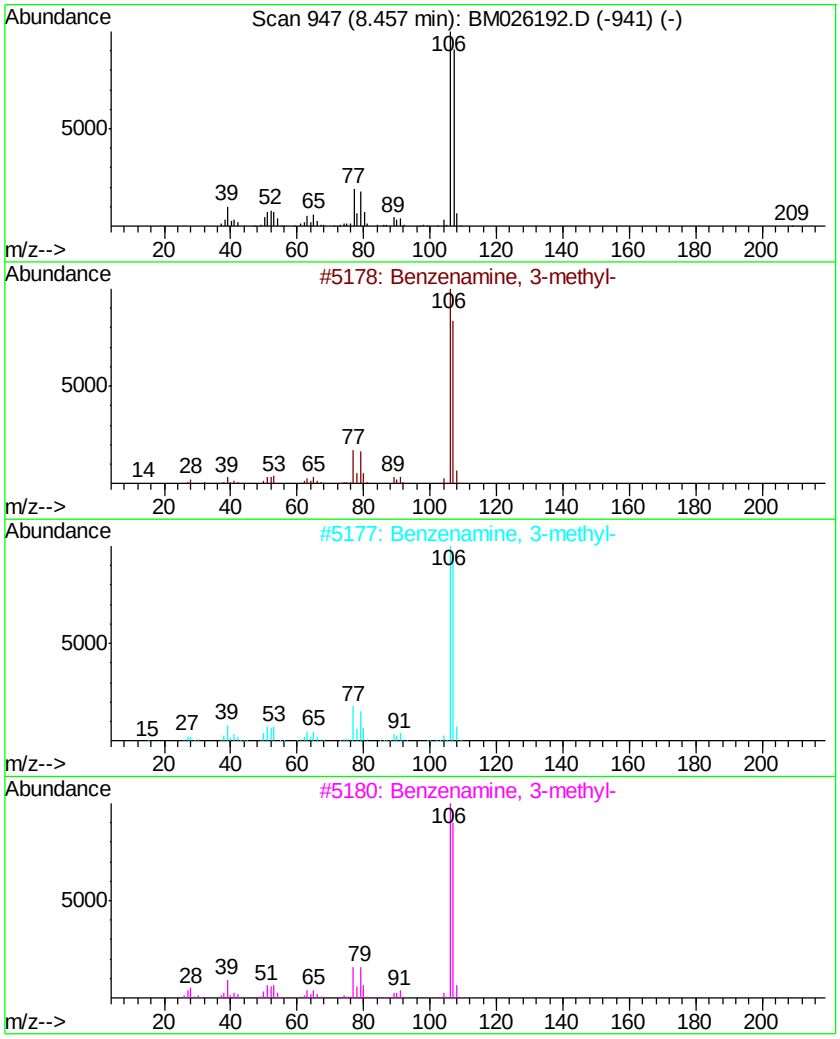
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

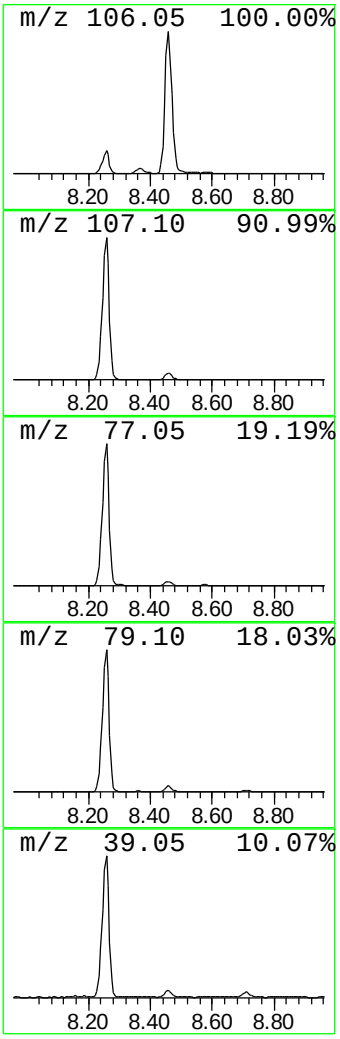
****

Peak Number 5 Benzenamine, 3-methyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.46	7.06 ng/ul	409117	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	97
4	Benzenamine, 3-methyl-	107	C7H9N	000108-44-1	96
5	o-Toluidine	107	C7H9N	000095-53-4	94





Library Search Compound Report

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Data File : BM026192.D
Acq On : 30 May 2020 03:47
Operator : MA/SJ
Sample : L2820-13
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : CB640

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

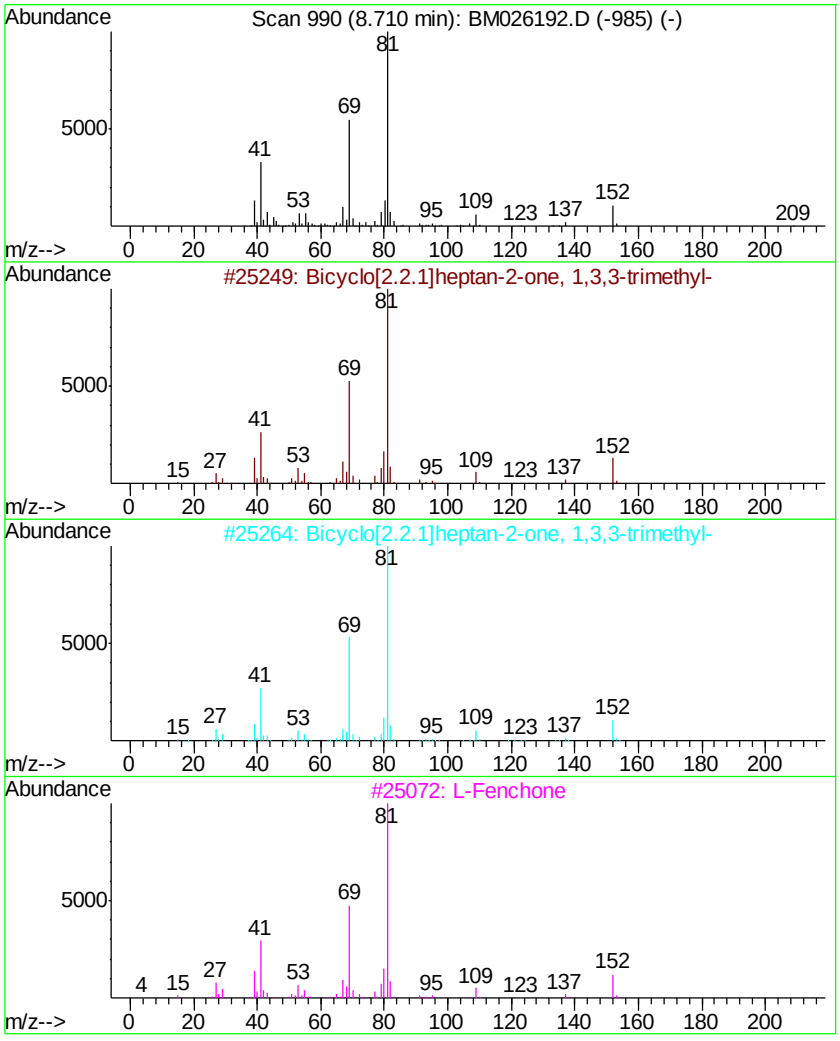
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

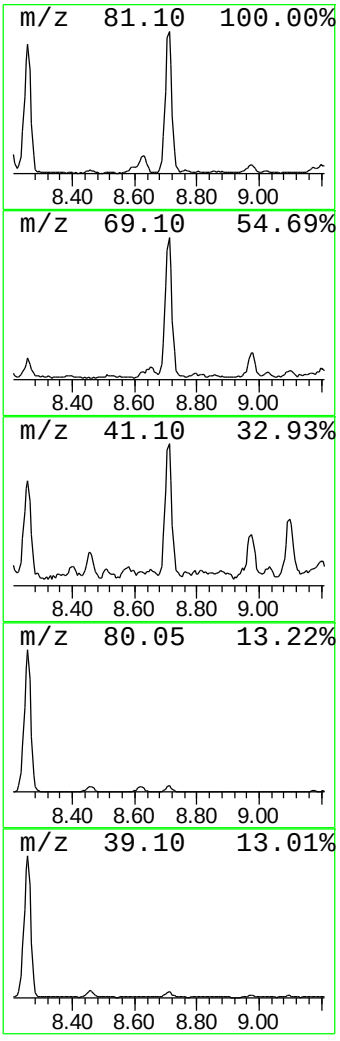
****

Peak Number 6 Bicyclo[2.2.1]heptan-2-one,... Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.71	4.05 ng/ul	234725	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	94
3	L-Fenchone	152	C10H16O	007787-20-4	91
4	D-Fenchone	152	C10H16O	004695-62-9	90
5	L-Fenchone	152	C10H16O	007787-20-4	90





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Sample : L2820-13
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : CB640

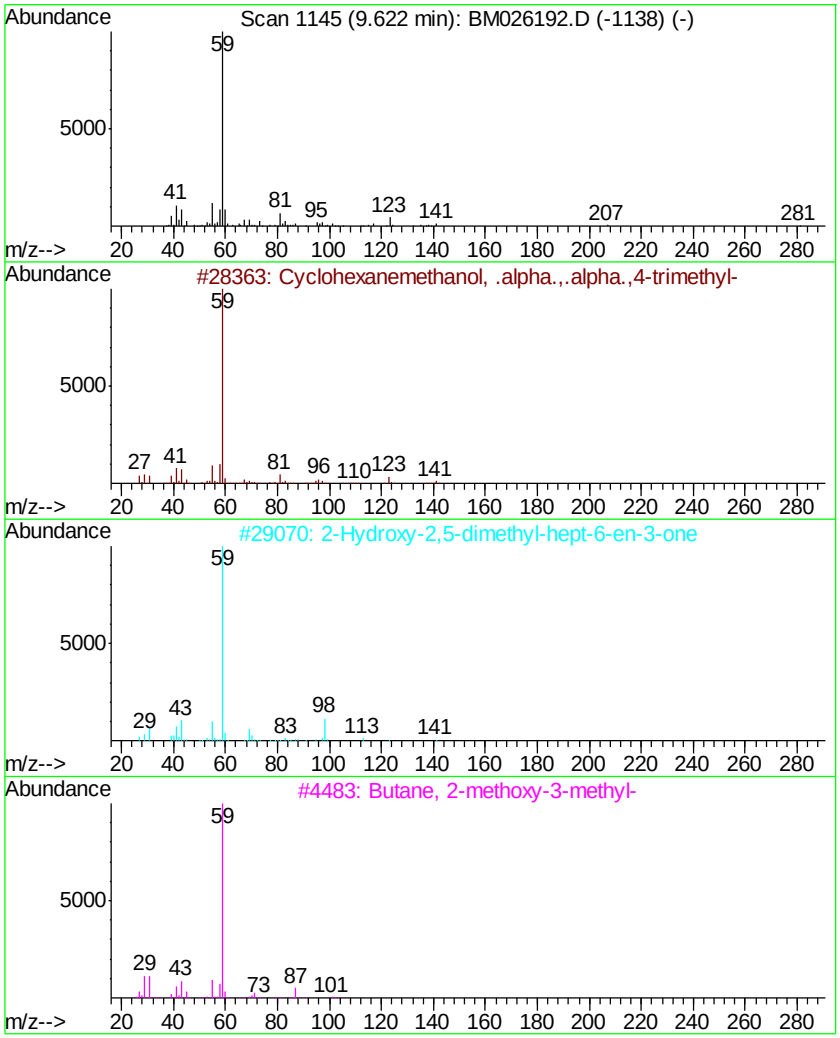
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM051320MA.M
Quant Title : SVOA CALIBRATION

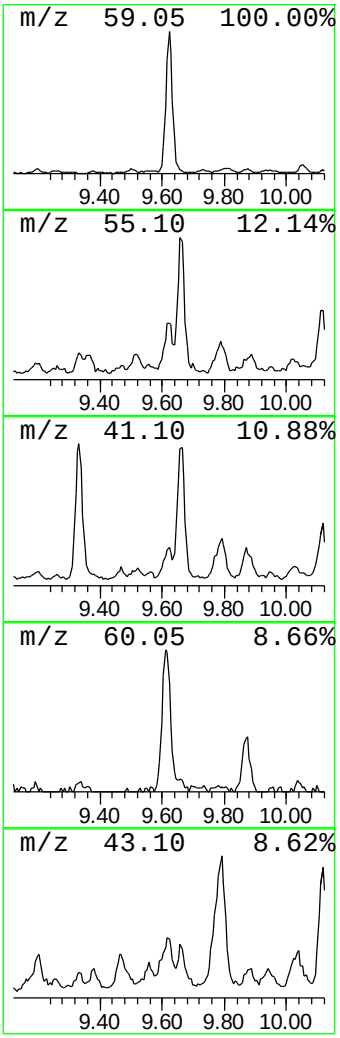
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 7 Cyclohexanemethanol, .alpha... Concentration Rank 23

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.62	2.52 ng/ul	190000	Naphthalene-d8	10.23

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclohexanemethanol, .alpha.,.al...	156	C10H20O	000498-81-7	72
2	2-Hydroxy-2,5-dimethyl-hept-6-en...	156	C9H16O2	1000192-55-8	64
3	Butane, 2-methoxy-3-methyl-	102	C6H14O	062016-49-3	59
4	Cyclohexanemethanol, .alpha.,.al...	156	C10H20O	005114-00-1	56
5	Undecanamide, 11-bromo-	263	C11H22BrNO	005875-26-3	39





Library Search Compound Report

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Data File : BM026192.D
Acq On : 30 May 2020 03:47
Operator : MA/SJ
Sample : L2820-13
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : CB640

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

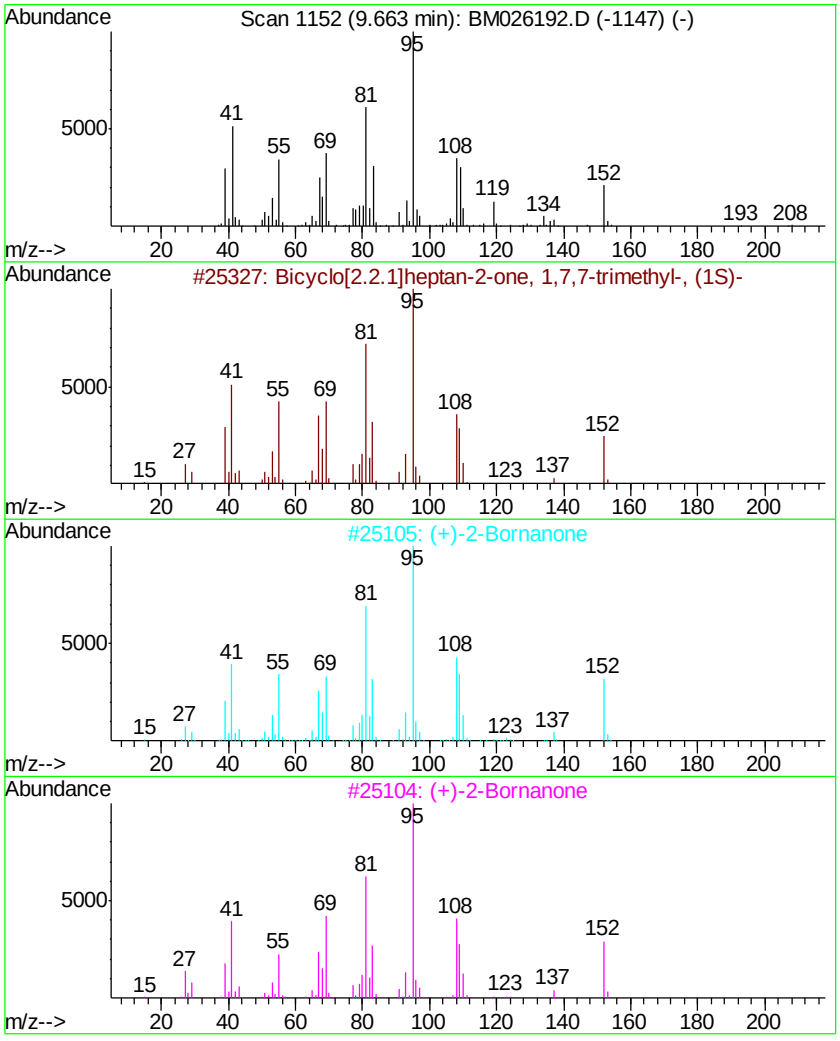
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

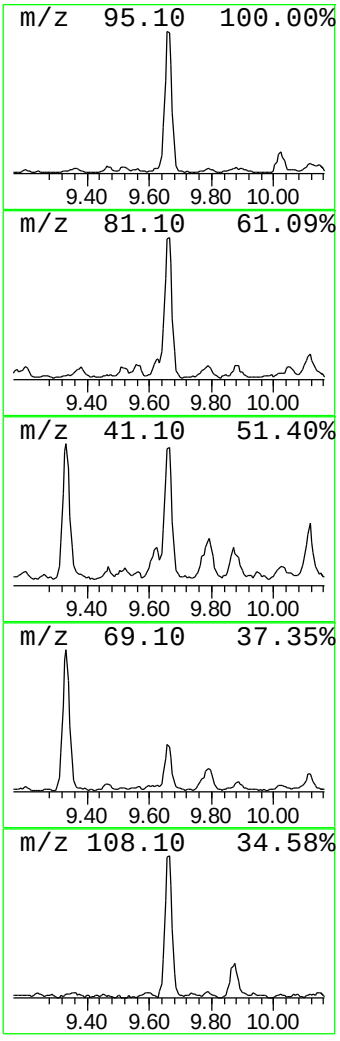
****

Peak Number 8 Bicyclo[2.2.1]heptan-2-one,... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.66	7.64 ng/ul	576580	Naphthalene-d8	10.23

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Bicyclo[2.2.1]heptan-2-one, 1,7,...	152	C10H16O	000464-48-2	97
2	(+)-2-Bornanone	152	C10H16O	000464-49-3	96
3	(+)-2-Bornanone	152	C10H16O	000464-49-3	96
4	Camphor	152	C10H16O	000076-22-2	96
5	(+)-2-Bornanone	152	C10H16O	000464-49-3	94





Library Search Compound Report

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Acq On : 30 May 2020 03:47
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Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : CB640

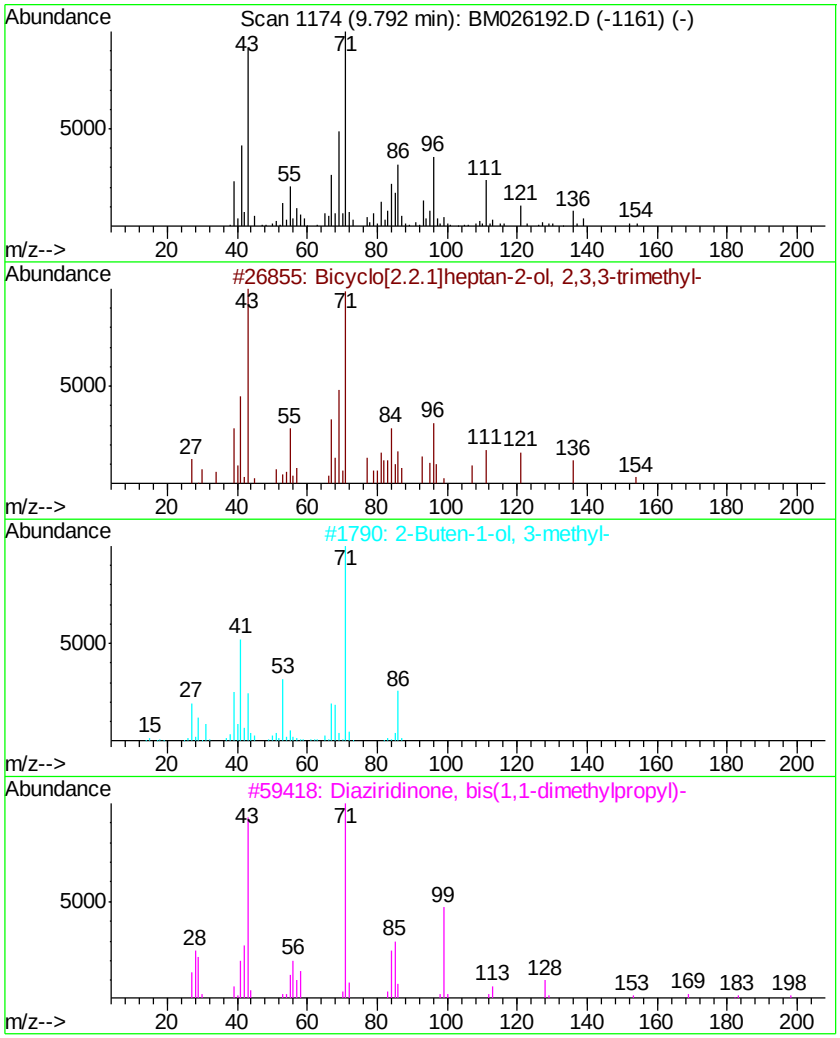
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM051320MA.M
Quant Title : SVOA CALIBRATION

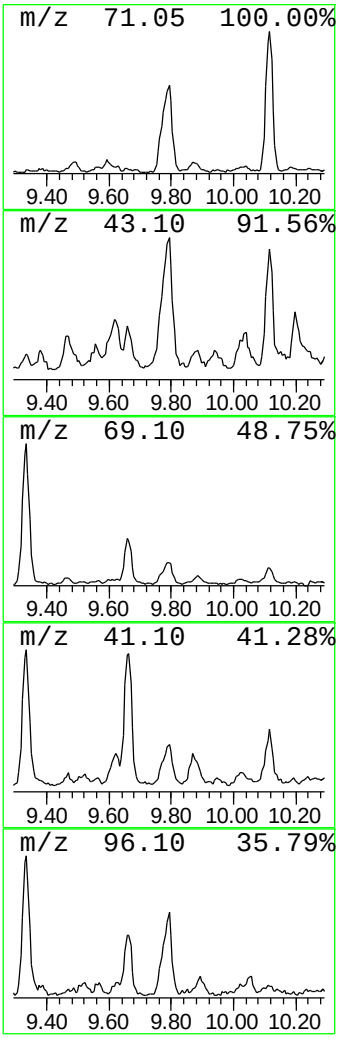
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 9 Bicyclo[2.2.1]heptan-2-ol, ... Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.79	4.09 ng/ul	308397	Naphthalene-d8	10.23

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Bicyclo[2.2.1]heptan-2-ol, 2,3,3...	154	C10H18O	000465-31-6	72
2	2-Buten-1-ol, 3-methyl-	86	C5H10O	000556-82-1	47
3	Diaziridinone, bis(1,1-dimethylp...	198	C11H22N2O	019656-75-8	43
4	2-Butene, 1-methoxy-, (E)-	86	C5H10O	010034-14-7	38
5	Acetic acid, (1,2-dimethyl-1-pro...	128	C7H12O2	003814-41-3	38





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Sample : L2820-13
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : CB640

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

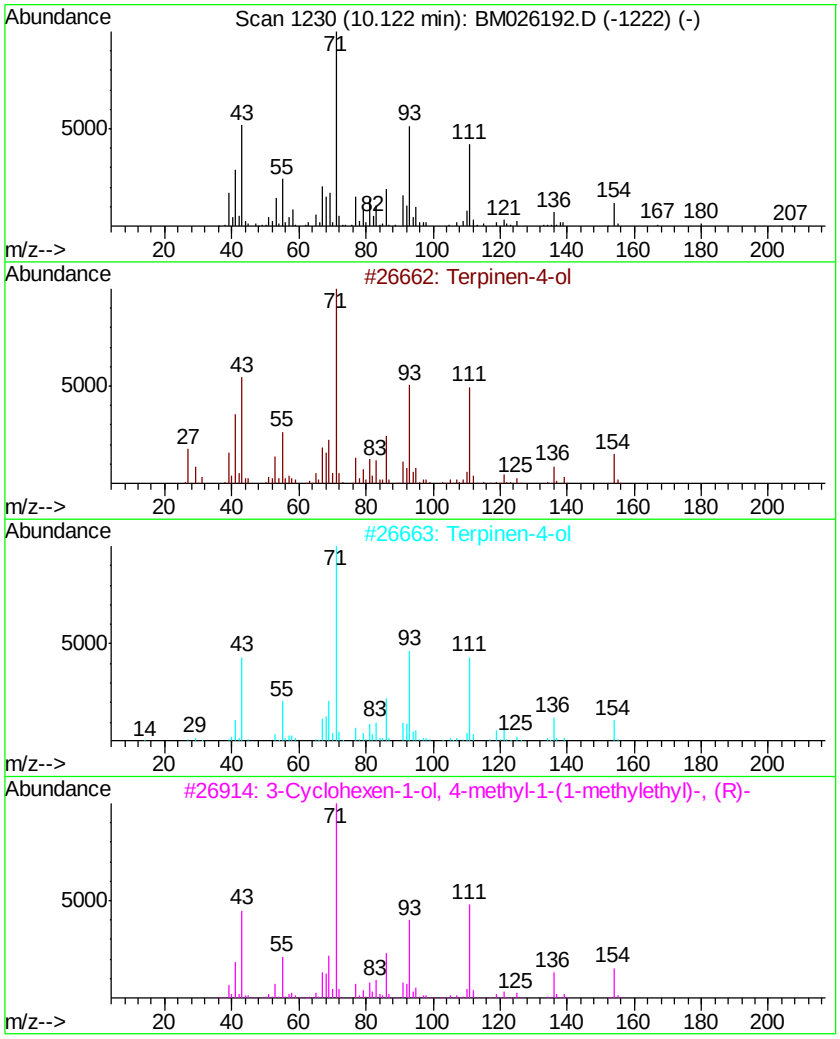
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

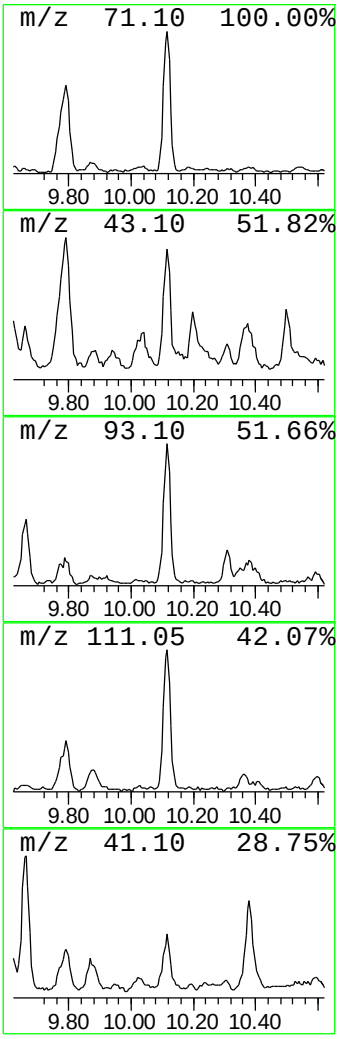
****

Peak Number 10 Terpinen-4-ol Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.12	4.08 ng/ul	307922	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	90
3	3-Cyclohexen-1-ol, 4-methyl-1-(1...	154	C10H18O	020126-76-5	87
4	3-Cyclohexen-1-ol, 4-methyl-1-(1...	154	C10H18O	020126-76-5	81
5	Terpinen-4-ol	154	C10H18O	000562-74-3	81





Library Search Compound Report

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Sample : L2820-13
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : CB640

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Quant Title : SVOA CALIBRATION

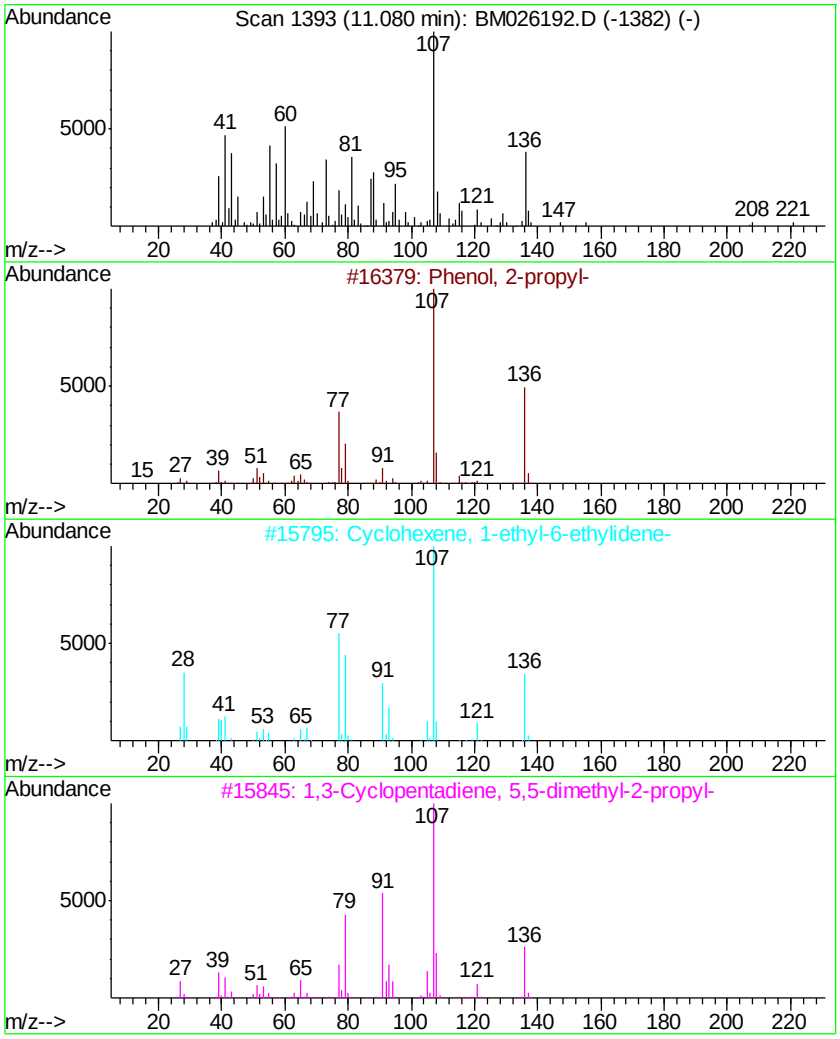
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

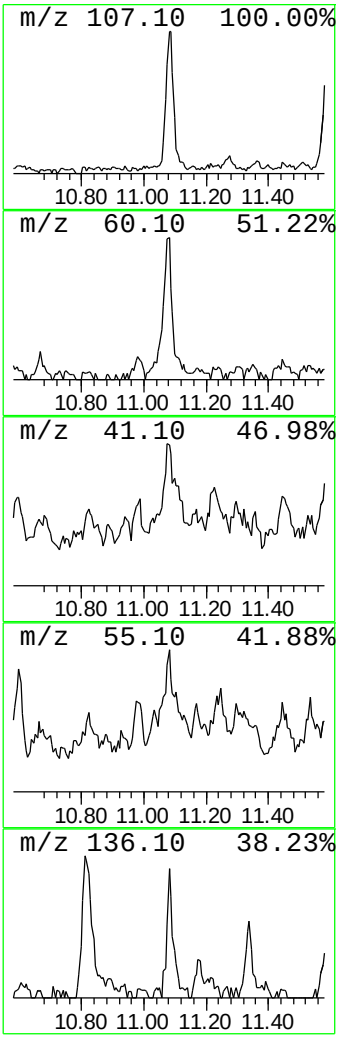
****

Peak Number 11 unknown-01 Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.08	2.78 ng/ul	209835	Naphthalene-d8	10.23

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, 2-propyl-	136	C9H12O	000644-35-9	45
2	Cyclohexene, 1-ethyl-6-ethylidene-	136	C10H16	061141-57-9	38
3	1,3-Cyclopentadiene, 5,5-dimethy...	136	C10H16	1000163-57-0	38
4	Phenol, 2-propyl-	136	C9H12O	000644-35-9	35
5	Pyridine, 2,6-dimethyl-	107	C7H9N	000108-48-5	22





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Sample : L2820-13
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : CB640

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

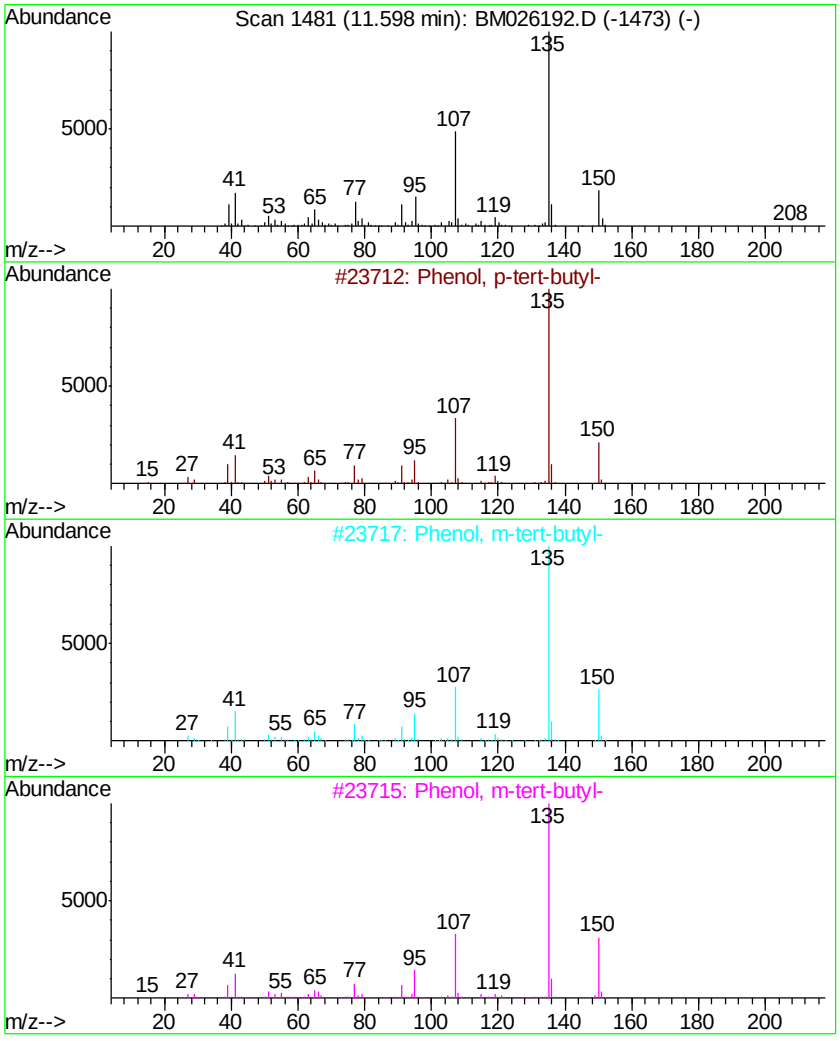
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

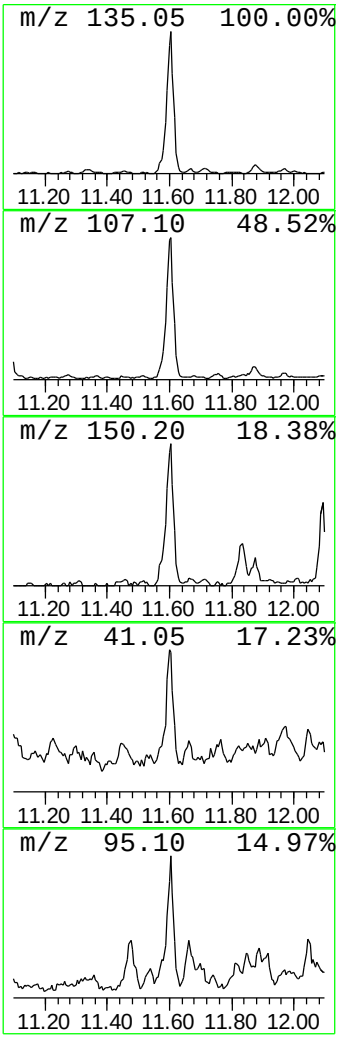
****

Peak Number 12 Phenol, p-tert-butyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.60	4.92 ng/ul	371324	Naphthalene-d8	10.23

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, p-tert-butyl-	150	C10H14O	000098-54-4	94
2	Phenol, m-tert-butyl-	150	C10H14O	000585-34-2	93
3	Phenol, m-tert-butyl-	150	C10H14O	000585-34-2	91
4	Phenol, p-tert-butyl-	150	C10H14O	000098-54-4	87
5	Phenol, p-tert-butyl-	150	C10H14O	000098-54-4	74





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Operator : MA/SJ
Sample : L2820-13
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : CB640

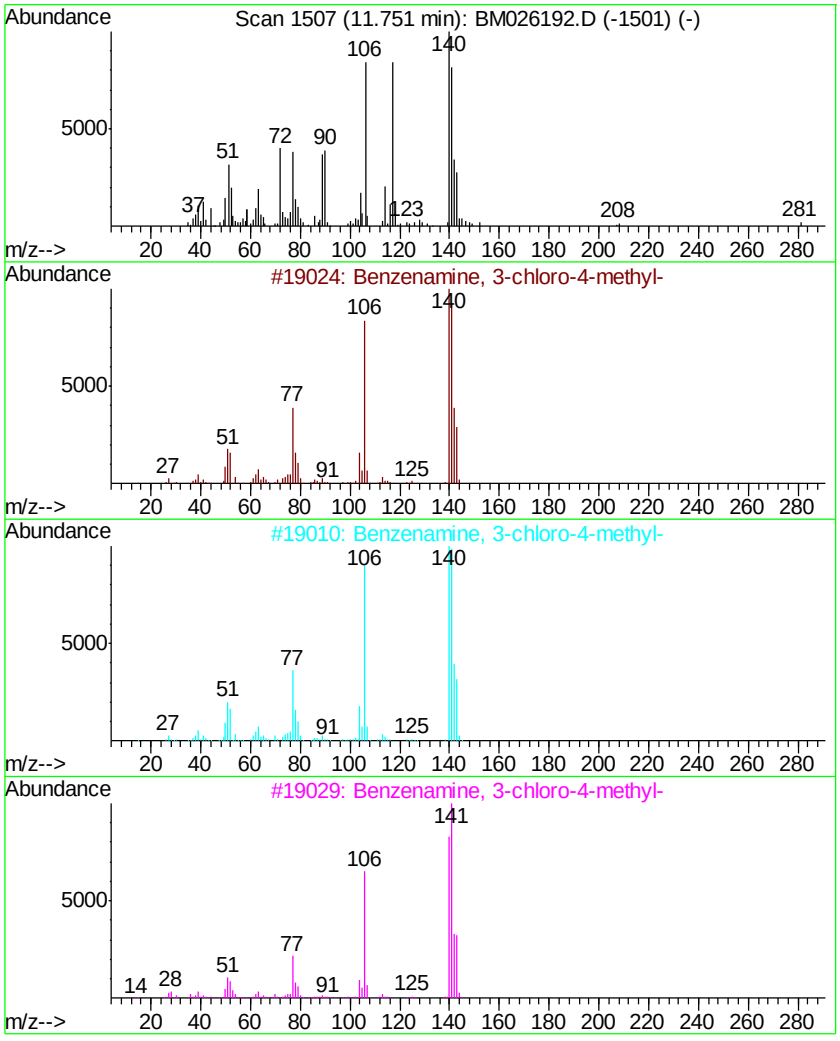
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM051320MA.M
Quant Title : SVOA CALIBRATION

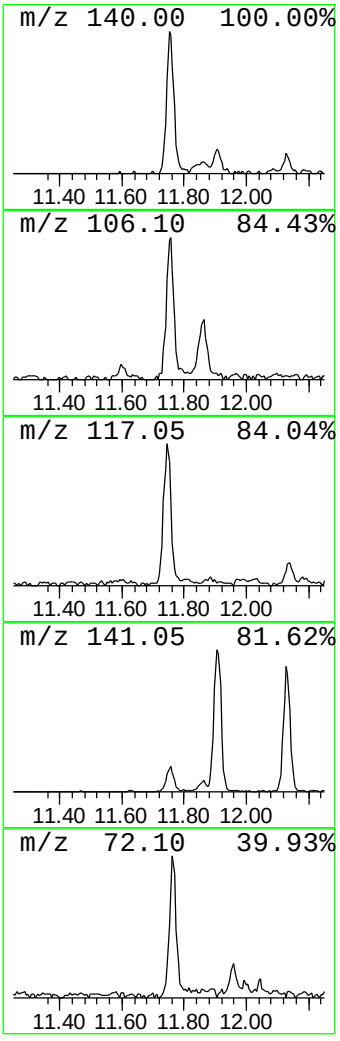
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 13 Benzenamine, 3-chloro-4-met... Concentration Rank 25

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.75	2.27 ng/ul	171042	Naphthalene-d8	10.23

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzenamine, 3-chloro-4-methyl-	141	C7H8ClN	000095-74-9	96
2	Benzenamine, 3-chloro-4-methyl-	141	C7H8ClN	000095-74-9	96
3	Benzenamine, 3-chloro-4-methyl-	141	C7H8ClN	000095-74-9	64
4	m-Toluidine, 4-chloro-	141	C7H8ClN	007149-75-9	64
5	Benzenamine, 2-chloro-4-methyl-	141	C7H8ClN	000615-65-6	64





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Operator : MA/SJ
Sample : L2820-13
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : CB640

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

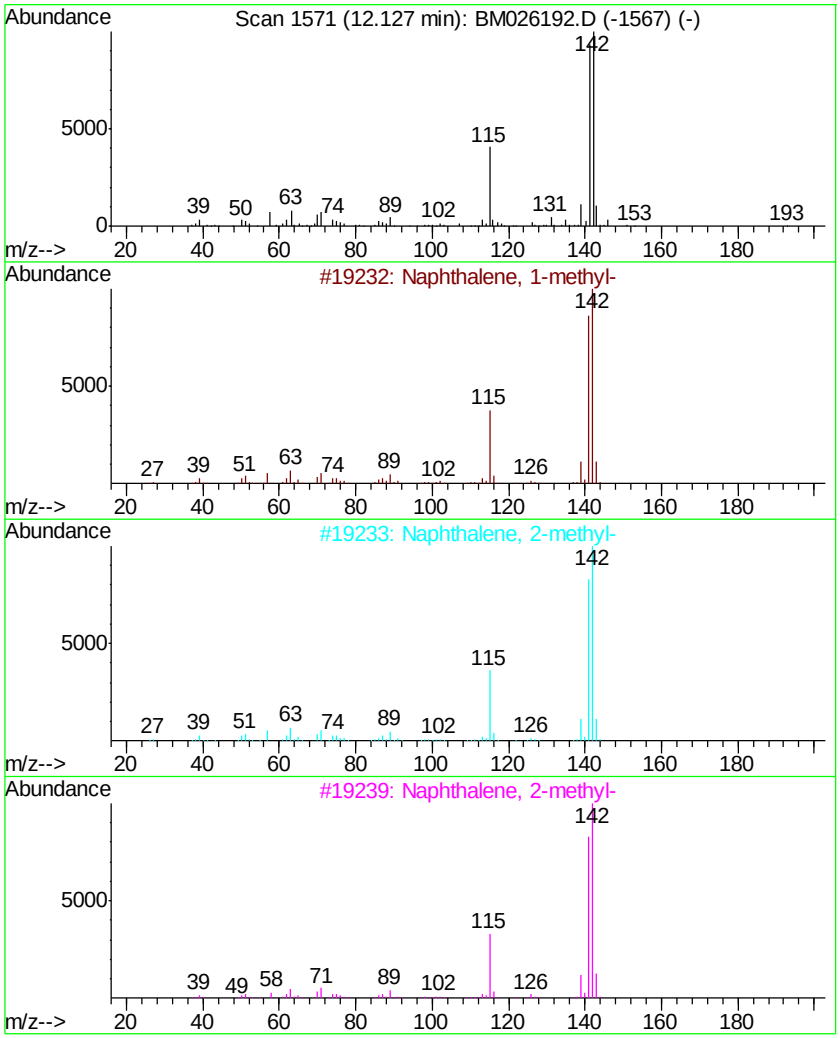
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

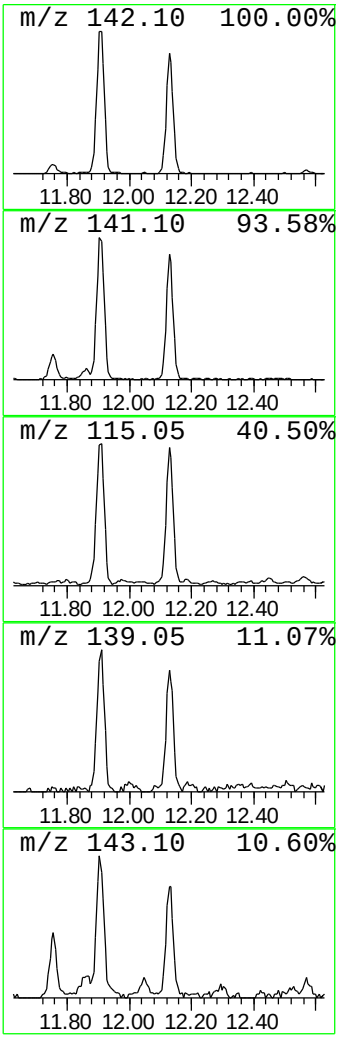
****

Peak Number 14 Naphthalene, 1-methyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.13	4.62 ng/ul	348461	Naphthalene-d8	10.23

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 1-methyl-	142	C11H10	000090-12-0	96
2	Naphthalene, 2-methyl-	142	C11H10	000091-57-6	96
3	Naphthalene, 2-methyl-	142	C11H10	000091-57-6	96
4	Naphthalene, 1-methyl-	142	C11H10	000090-12-0	91
5	Benzocycloheptatriene	142	C11H10	000264-09-5	87





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Sample : L2820-13
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : CB640

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

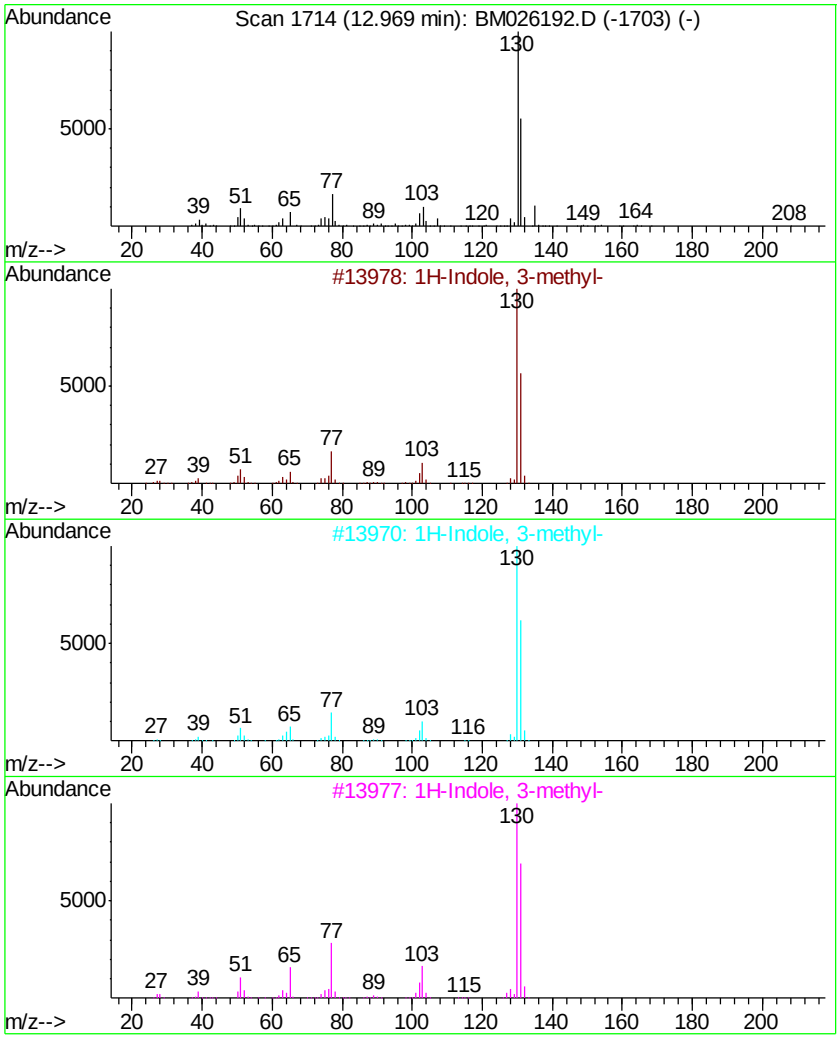
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

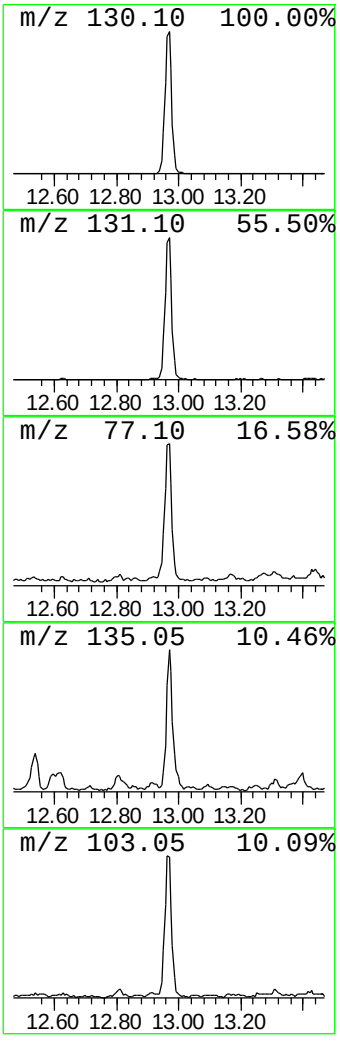
****

Peak Number 15 1H-Indole, 3-methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.97	6.55 ng/ul	1010690	Acenaphthene-d10	14.12

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Indole, 3-methyl-	131	C9H9N	000083-34-1	94
2	1H-Indole, 3-methyl-	131	C9H9N	000083-34-1	94
3	1H-Indole, 3-methyl-	131	C9H9N	000083-34-1	90
4	1H-Indole, 2-methyl-	131	C9H9N	000095-20-5	90
5	1H-Indole, 3-methyl-	131	C9H9N	000083-34-1	83





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ClientSampleId : CB640

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Quant Title : SVOA CALIBRATION

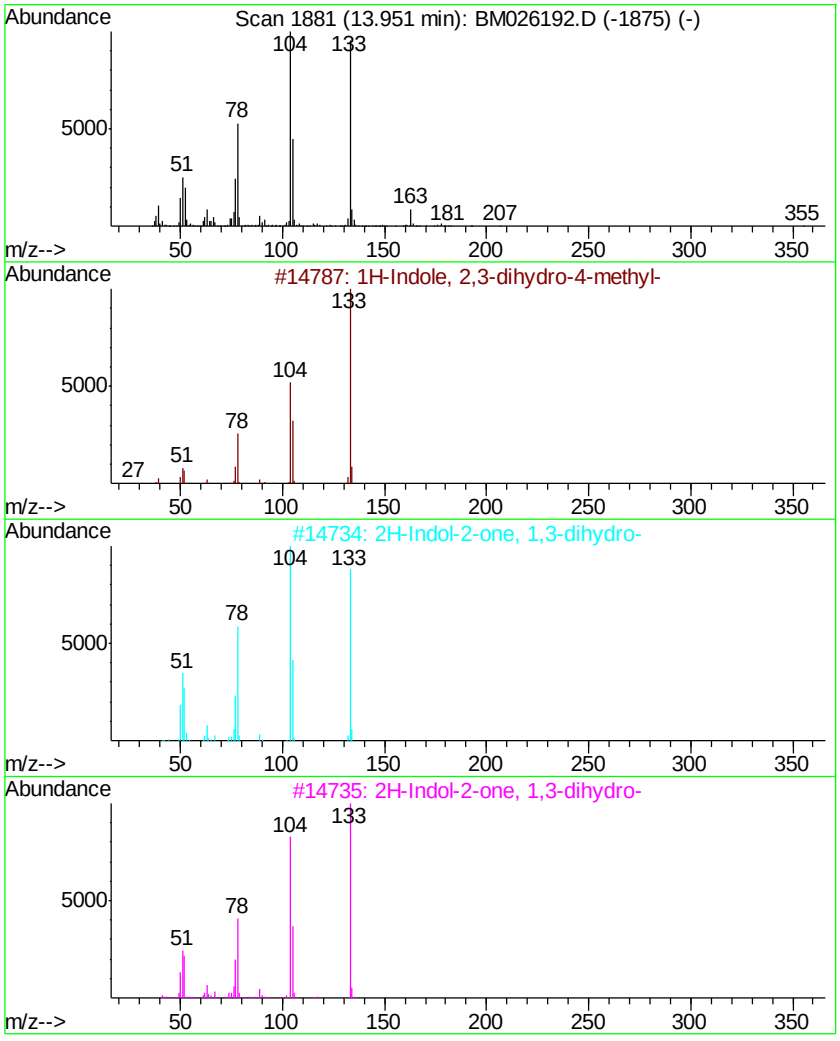
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

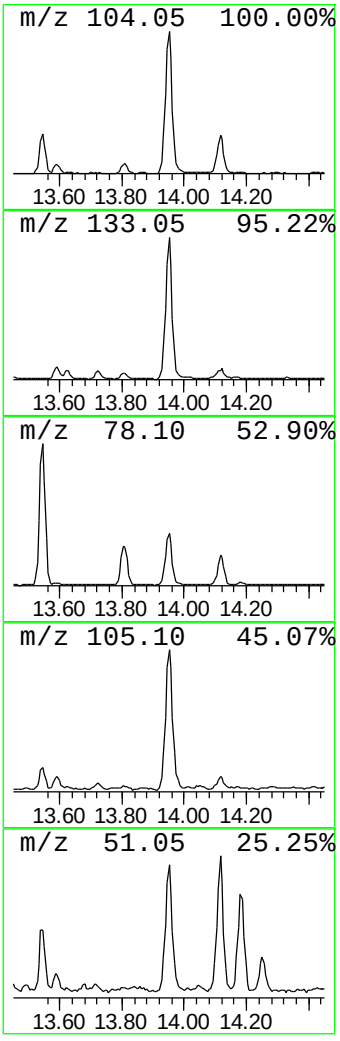
****

Peak Number 16 1H-Indole, 2,3-dihydro-4-me... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.95	5.57 ng/ul	860011	Acenaphthene-d10	14.12

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Indole, 2,3-dihydro-4-methyl-	133	C9H11N	062108-16-1	97
2	2H-Indol-2-one, 1,3-dihydro-	133	C8H7NO	000059-48-3	97
3	2H-Indol-2-one, 1,3-dihydro-	133	C8H7NO	000059-48-3	95
4	2H-Indol-2-one, 1,3-dihydro-	133	C8H7NO	000059-48-3	94
5	1H-Benzotriazole, 5-methyl-	133	C7H7N3	000136-85-6	91





Library Search Compound Report

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Data File : BM026192.D
Acq On : 30 May 2020 03:47
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Sample : L2820-13
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : CB640

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

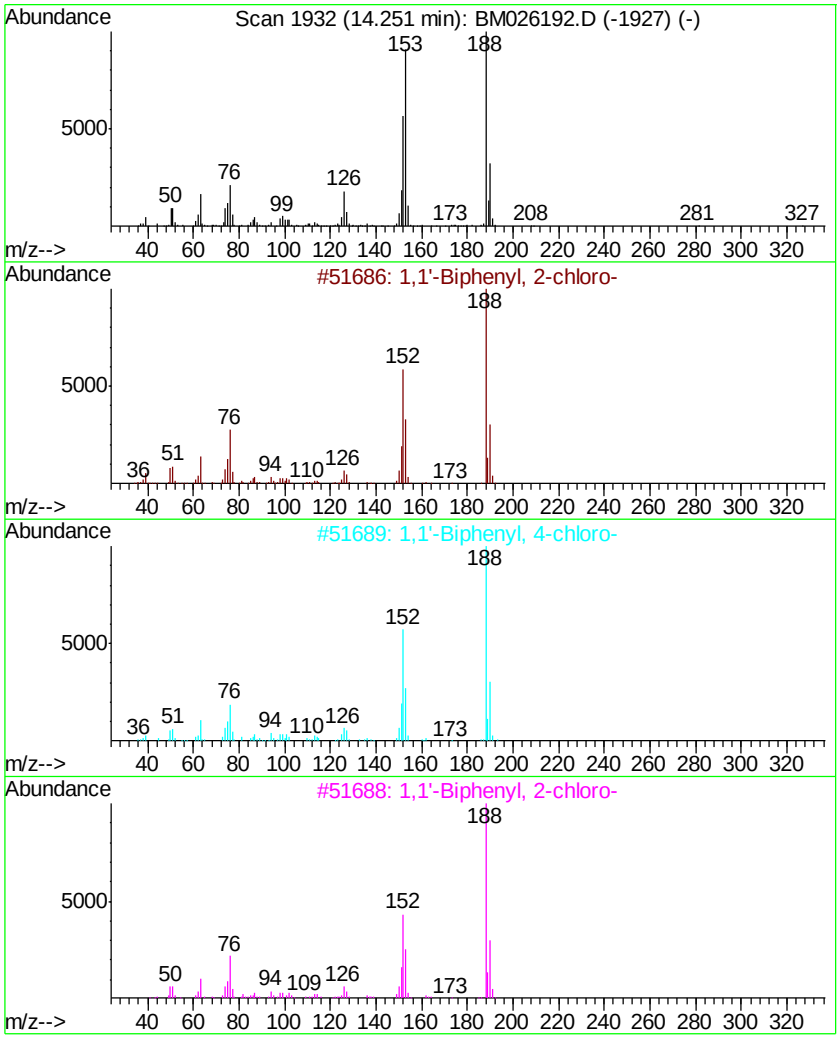
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

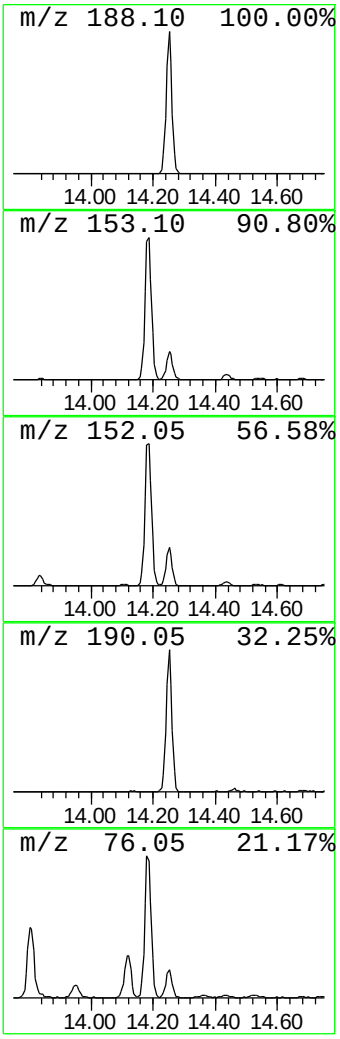
****

Peak Number 17 1,1'-Biphenyl, 2-chloro- Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.25	3.28 ng/ul	506026	Acenaphthene-d10	14.12

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl, 2-chloro-	188	C12H9Cl	002051-60-7	89
2	1,1'-Biphenyl, 4-chloro-	188	C12H9Cl	002051-62-9	86
3	1,1'-Biphenyl, 2-chloro-	188	C12H9Cl	002051-60-7	83
4	1,1'-Biphenyl, 2-chloro-	188	C12H9Cl	002051-60-7	83
5	3-Chlorobiphenyl	188	C12H9Cl	002051-61-8	83





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Acq On : 30 May 2020 03:47
Operator : MA/SJ
Sample : L2820-13
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : CB640

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

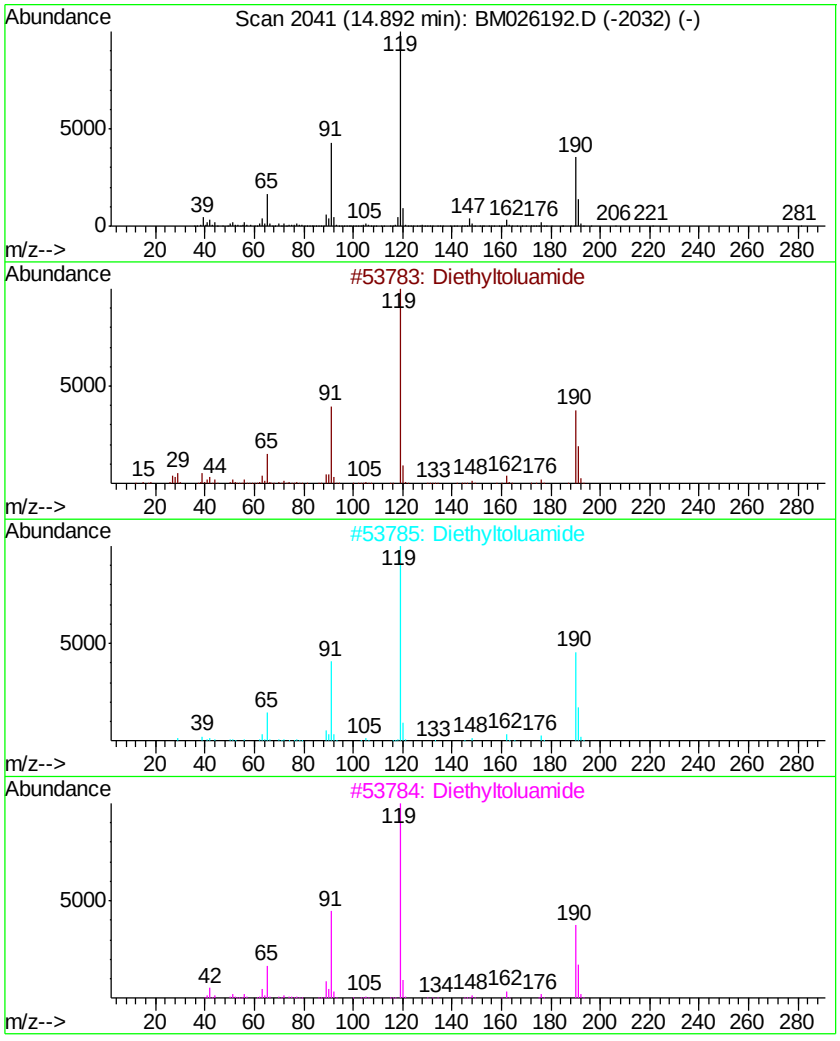
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

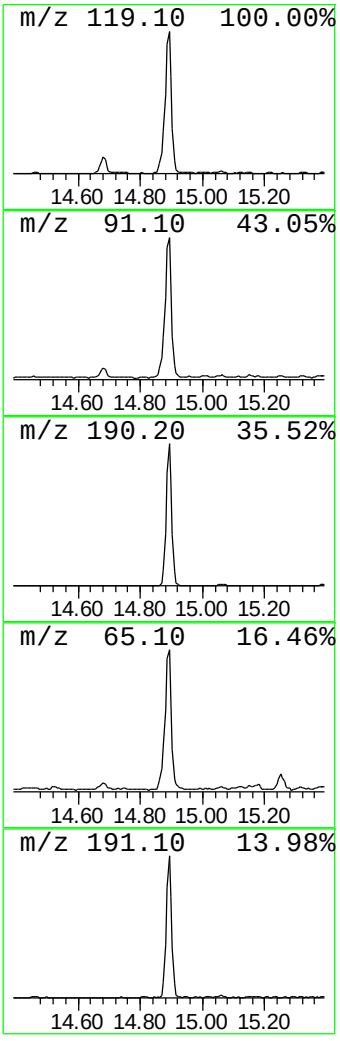
****

Peak Number 18 Diethyltoluamide Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.89	14.86 ng/ul	2293640	Acenaphthene-d10	14.12

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Diethyltoluamide	191	C12H17NO	000134-62-3	95
2	Diethyltoluamide	191	C12H17NO	000134-62-3	94
3	Diethyltoluamide	191	C12H17NO	000134-62-3	90
4	Benzamide, N,N-diethyl-4-methyl-	191	C12H17NO	002728-05-4	76
5	4'-Methylbutyrophenone	162	C11H14O	004160-52-5	58





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Acq On : 30 May 2020 03:47
Operator : MA/SJ
Sample : L2820-13
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : CB640

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

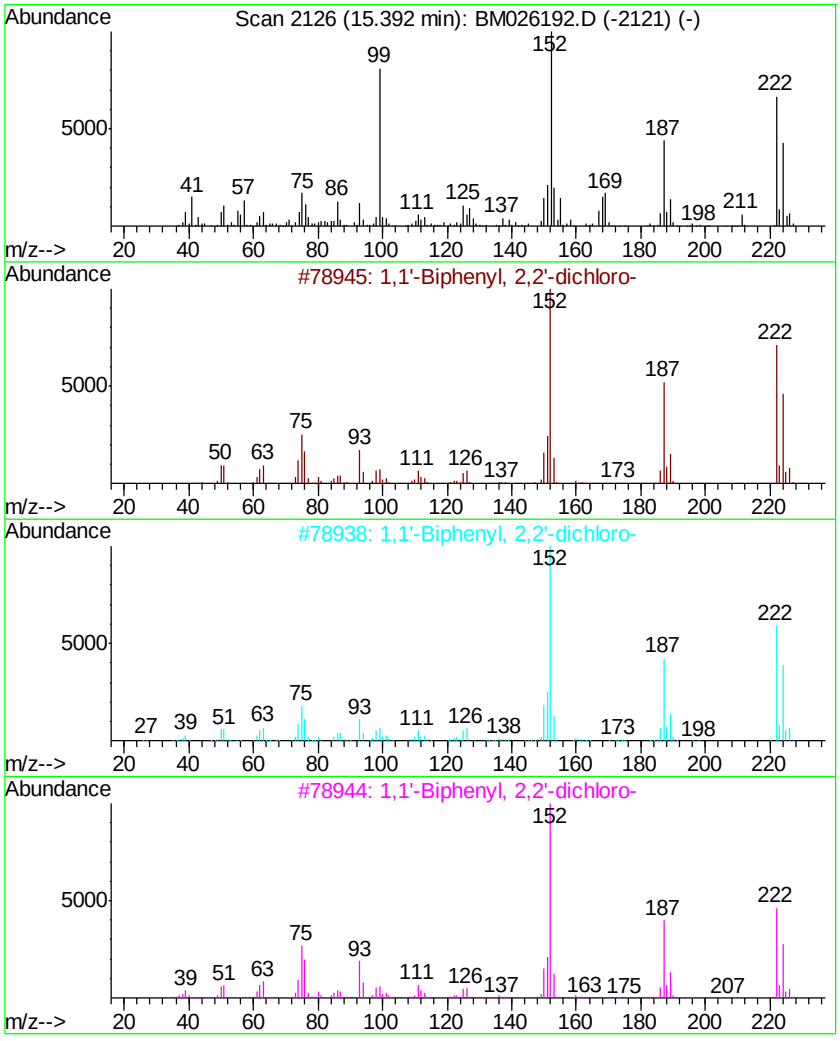
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

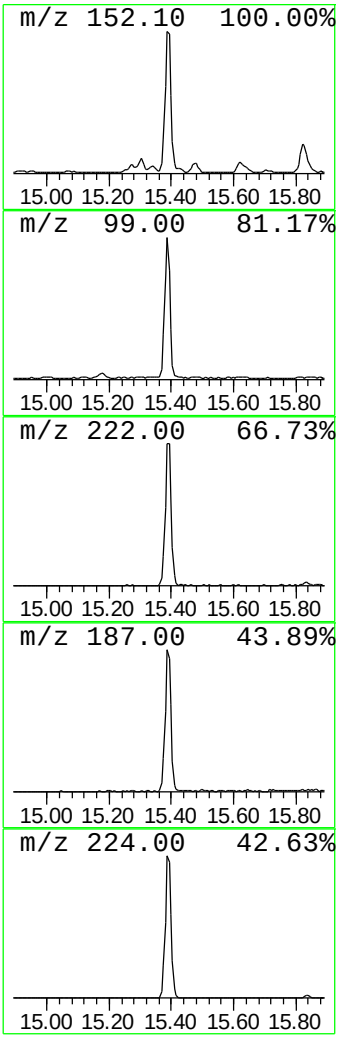
****

Peak Number 19 1,1'-Biphenyl, 2,2'-dichloro- Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.39	3.67 ng/ul	566686	Acenaphthene-d10	14.12

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl, 2,2'-dichloro-	222	C12H8Cl2	013029-08-8	99
2	1,1'-Biphenyl, 2,2'-dichloro-	222	C12H8Cl2	013029-08-8	99
3	1,1'-Biphenyl, 2,2'-dichloro-	222	C12H8Cl2	013029-08-8	98
4	1,1'-Biphenyl, 3,4'-dichloro-	222	C12H8Cl2	002974-90-5	95
5	1,1'-Biphenyl, 4,4'-dichloro-	222	C12H8Cl2	002050-68-2	95





Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM052920\
Data File : BM026192.D
Acq On : 30 May 2020 03:47
Operator : MA/SJ
Sample : L2820-13
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : CB640

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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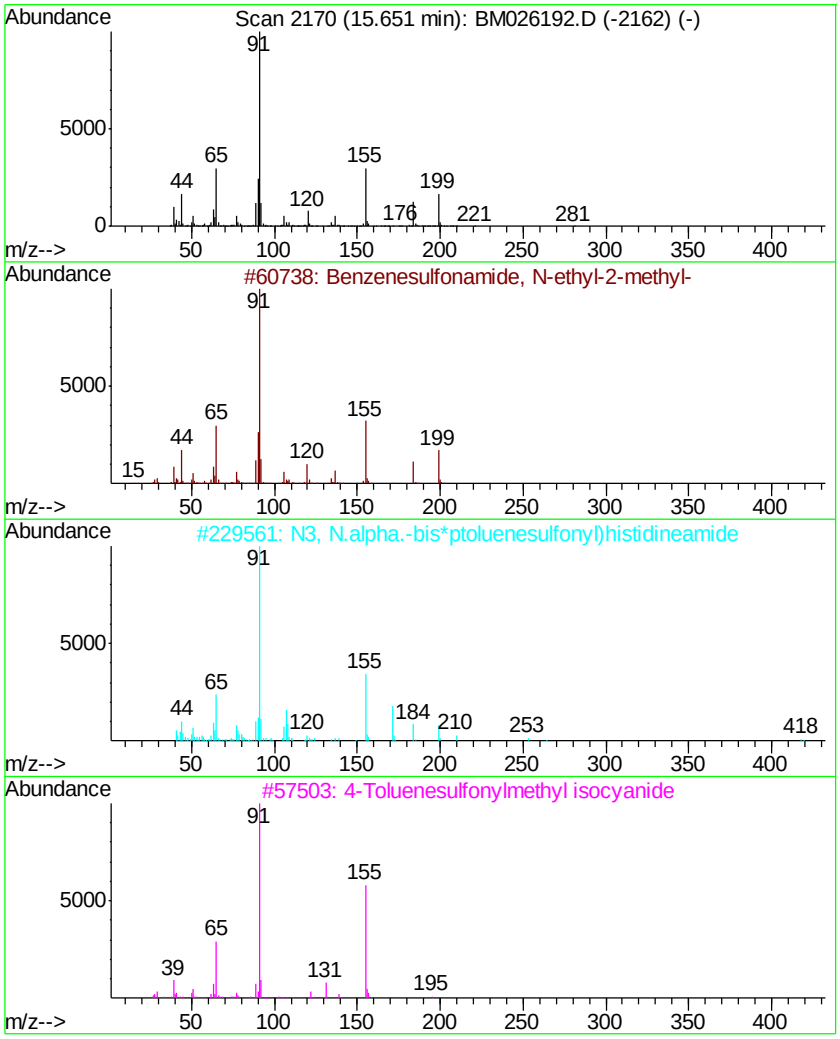
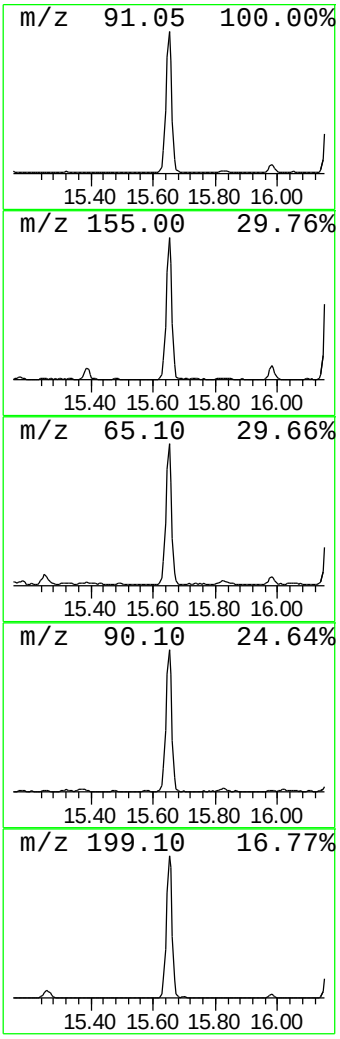
Peak Number 20 Benzenesulfonamide, N-ethyl... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.65	17.56 ng/ul	1821450	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	N3, N.alpha.-bis*ptoluenesulfony...	462	C20H22N4O5S2	1000130-21-9	53
3	4-Toluenesulfonylmethyl isocyanide	195	C9H9NO2S	036635-61-7	52
4	p-Toluenesulfonylacetonitrile	195	C9H9NO2S	005697-44-9	50
5	(O-p-Tosylisonitroso)malononitrile	249	C10H7N3O3S	020893-01-0	47





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Data File : BM026192.D
Acq On : 30 May 2020 03:47
Operator : MA/SJ
Sample : L2820-13
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : CB640

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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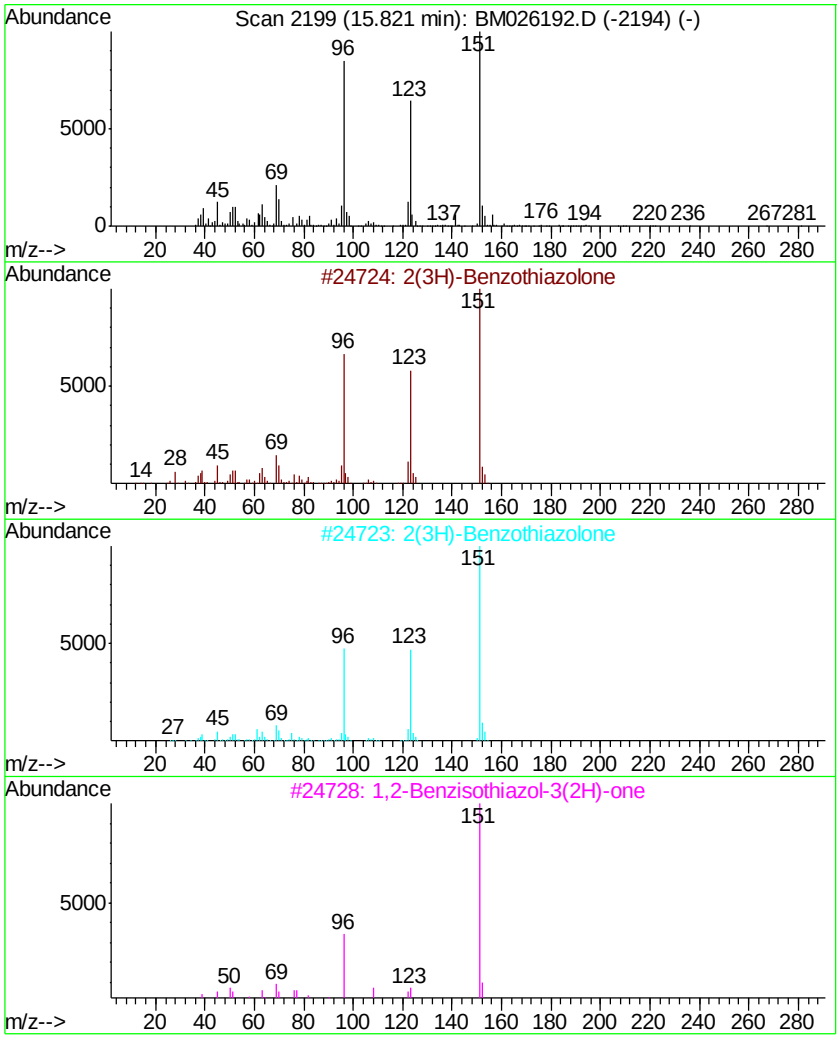
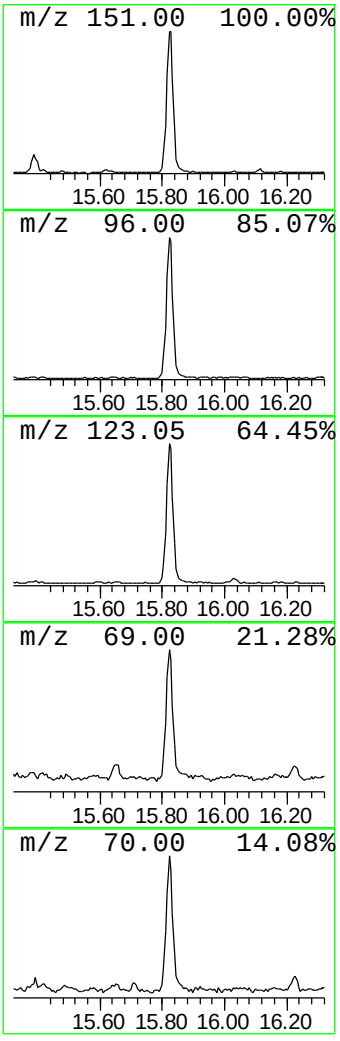
Peak Number 21 2(3H)-Benzothiazolone Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.82	8.87 ng/ul	919380	Phenanthrene-d10	16.87

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2(3H)-Benzothiazolone	151	C7H5NOS	000934-34-9	96
2	2(3H)-Benzothiazolone	151	C7H5NOS	000934-34-9	94
3	1,2-Benzisothiazol-3(2H)-one	151	C7H5NOS	002634-33-5	76
4	t-Butylhydroquinone	166	C10H14O2	001948-33-0	43
5	4-Methoxyformanilide	151	C8H9NO2	005470-34-8	35





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Data File : BM026192.D
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Operator : MA/SJ
Sample : L2820-13
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : CB640

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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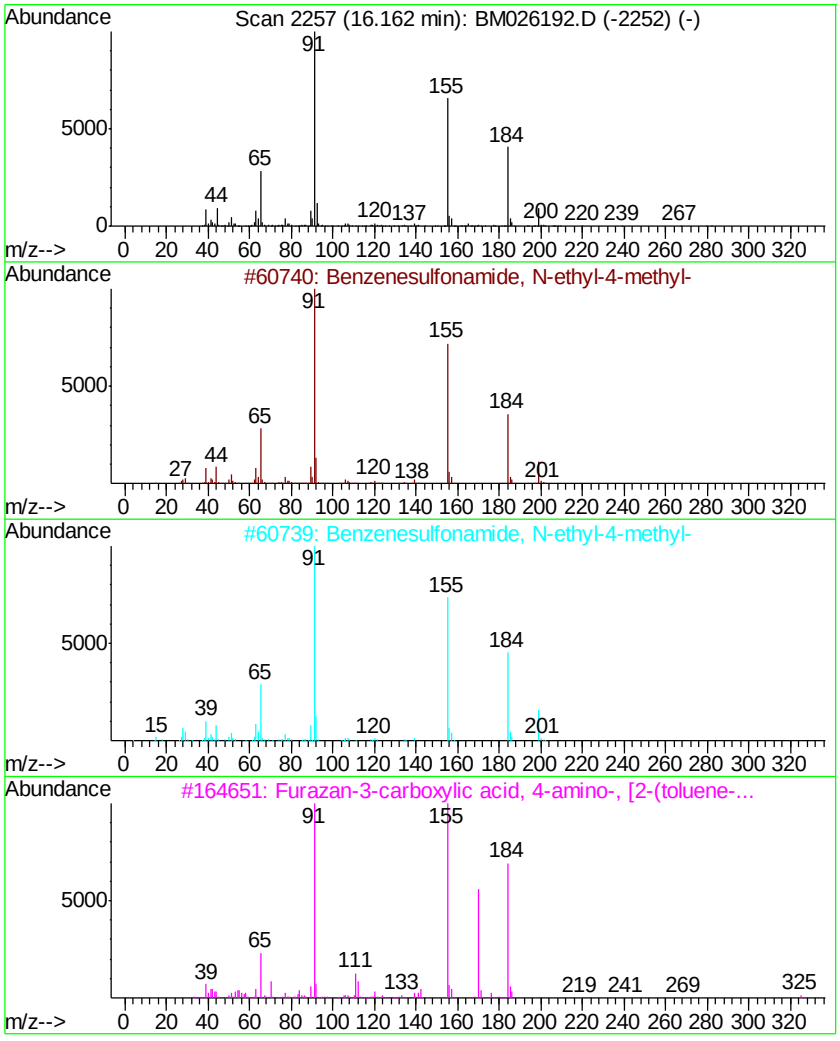
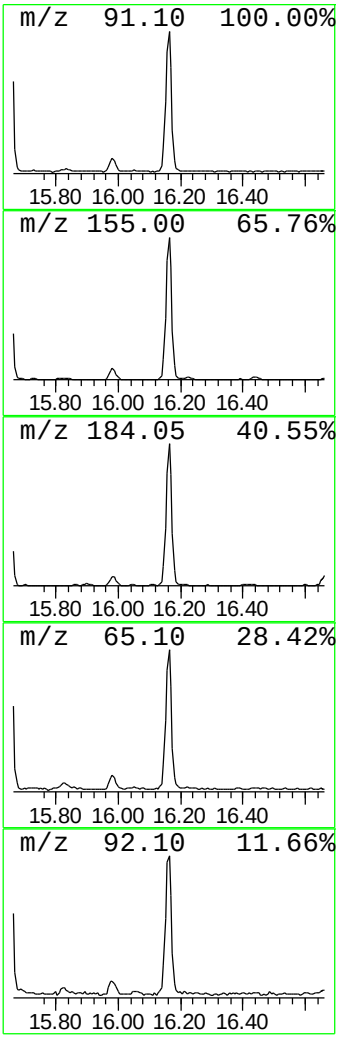
Peak Number 22 Benzenesulfonamide, N-ethyl... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.16	9.70 ng/ul	1005680	Phenanthrene-d10	16.87

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzenesulfonamide, N-ethyl-4-me...	199	C9H13NO2S	000080-39-7	97
2	Benzenesulfonamide, N-ethyl-4-me...	199	C9H13NO2S	000080-39-7	96
3	Furazan-3-carboxylic acid, 4-ami...	325	C12H15N5O4S	1000304-69-4	78
4	(Toluene-4-sulfonylamino)-acetic...	283	C13H17NO4S	1000190-48-0	64
5	4-Methyl-N-(2-oxobutyl)benzenesu...	241	C11H15NO3S	1000192-14-5	59





Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM052920\
Data File : BM026192.D
Acq On : 30 May 2020 03:47
Operator : MA/SJ
Sample : L2820-13
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : CB640

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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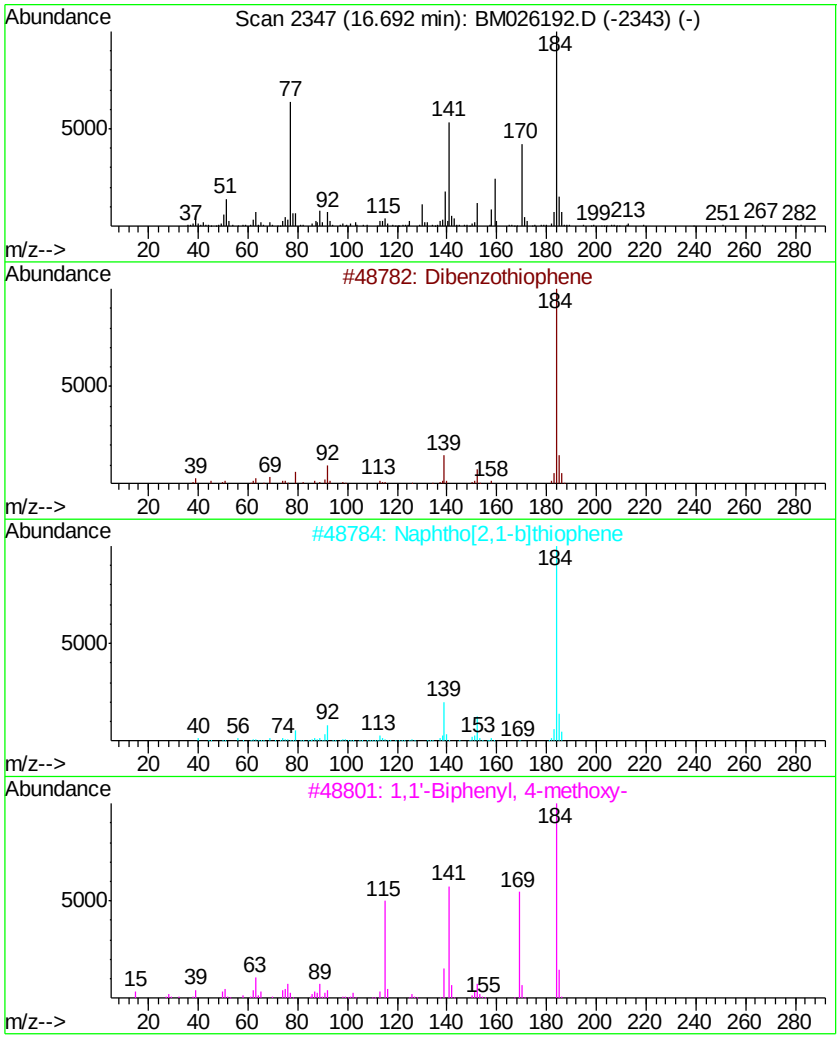
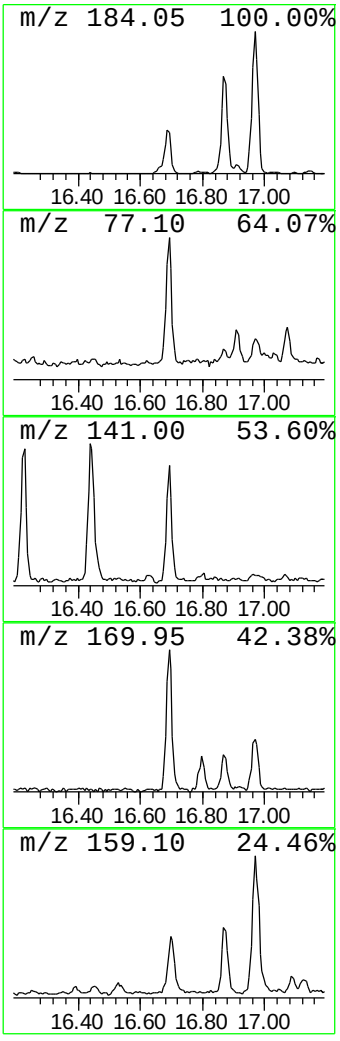
Peak Number 23 Dibenzothiophene Concentration Rank 24

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.69	2.48 ng/ul	256843	Phenanthrene-d10	16.87

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Dibenzothiophene	184	C12H8S	000132-65-0	80
2	Naphtho[2,1-b]thiophene	184	C12H8S	000233-02-3	55
3	1,1'-Biphenyl, 4-methoxy-	184	C13H12O	000613-37-6	46
4	Benzenesulfonamide, N-butyl-	213	C10H15NO2S	003622-84-2	46
5	Benzenesulfonamide, N-butyl-	213	C10H15NO2S	003622-84-2	46





Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM052920\
Data File : BM026192.D
Acq On : 30 May 2020 03:47
Operator : MA/SJ
Sample : L2820-13
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : CB640

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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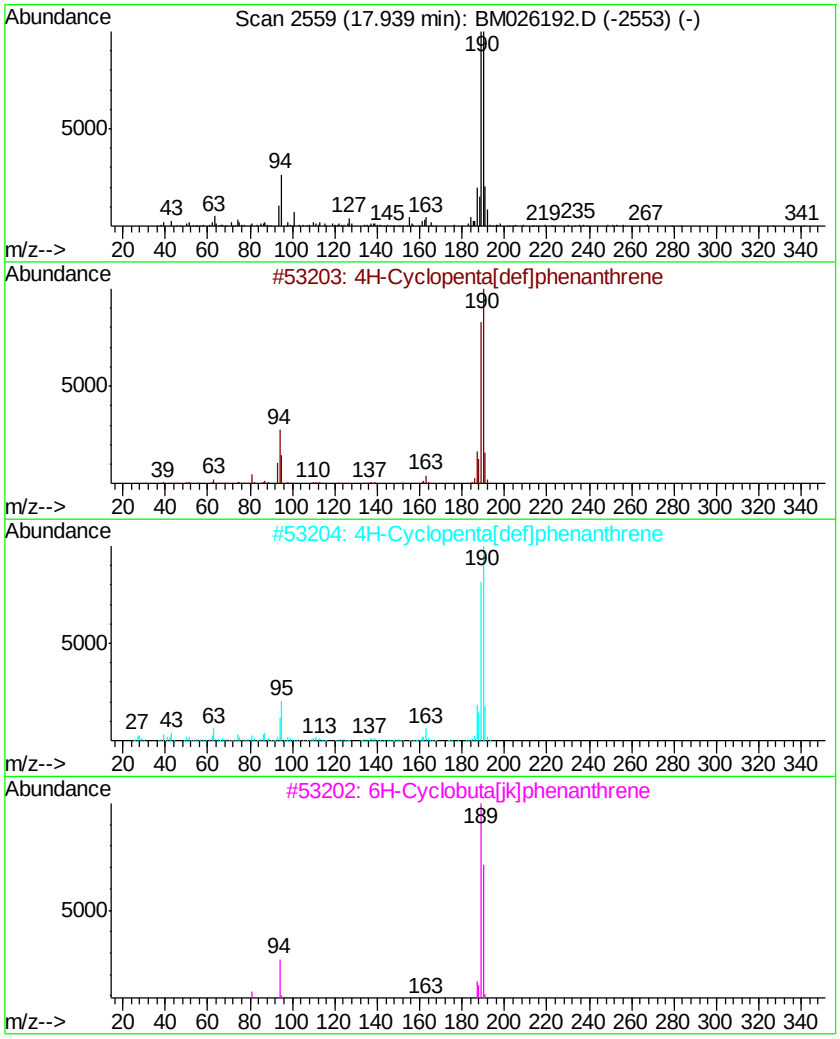
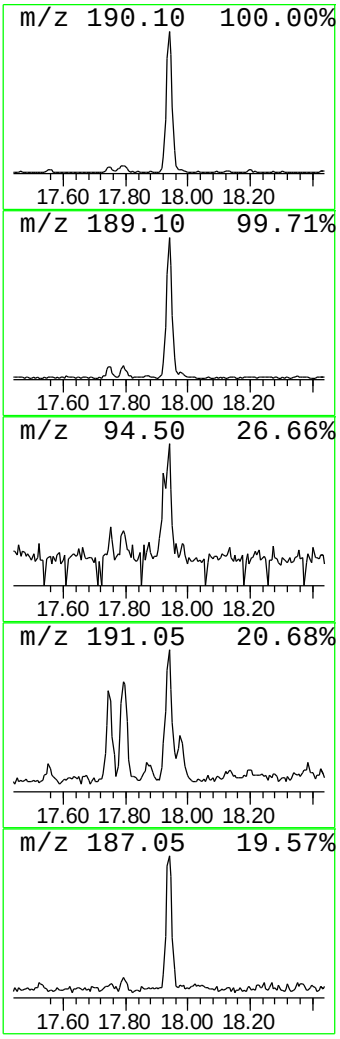
Peak Number 24 4H-Cyclopenta[def]phenanthrene Concentration Rank 22

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.94	2.58 ng/ul	268044	Phenanthrene-d10	16.87

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	91
2	4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	81
3	6H-Cyclobuta[jk]phenanthrene	190	C15H10	083469-43-6	78
4	4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	53
5	Phenol, 4-(2-thienylmethyl)-	190	C11H10OS	091680-55-6	45





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Operator : MA/SJ
Sample : L2820-13
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : CB640

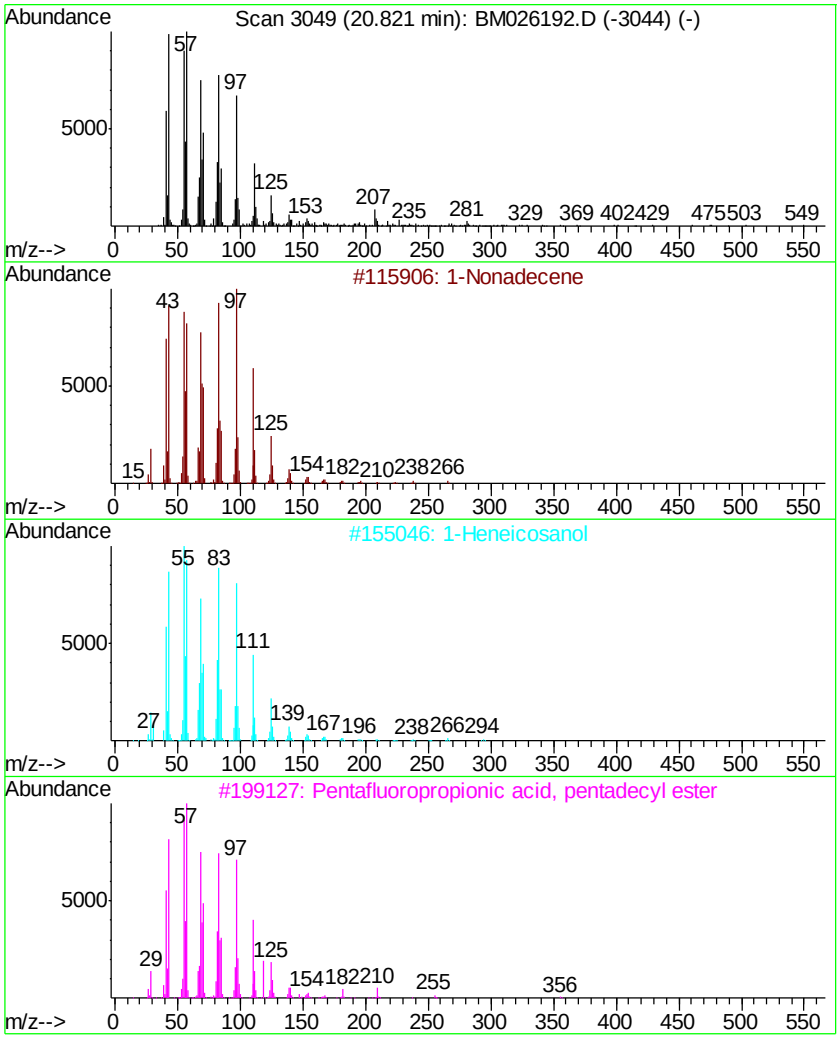
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM051320MA.M
Quant Title : SVOA CALIBRATION

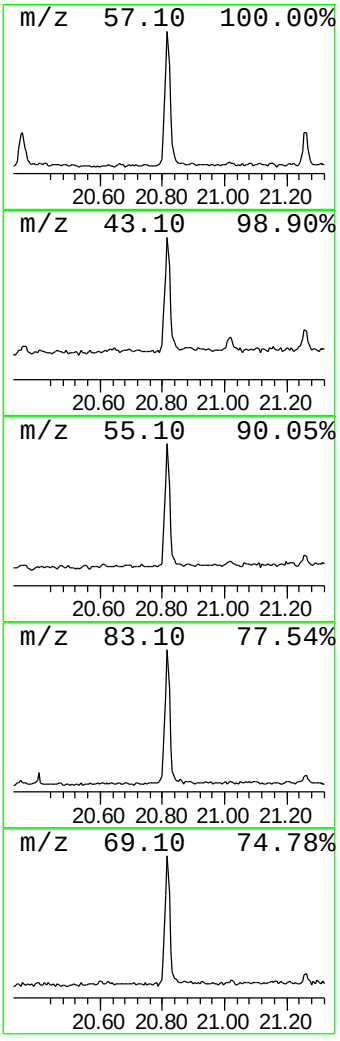
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 25 (DEL) Alkane: Straight-Chai... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.82	6.13 ng/ul	711784	Chrysene-d12	21.07

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Nonadecene	266	C19H38	018435-45-5	97
2	1-Heneicosanol	312	C21H44O	015594-90-8	94
3	Pentafluoropropionic acid, penta...	374	C18H31F5O2	959092-08-1	94
4	Cyclopentadecane	210	C15H30	000295-48-7	93
5	2- Chloropropionic acid, pentade...	318	C18H35ClO2	1000292-44-1	91





Tentatively Identified Compound (LSC) summary

Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM052920\
Data File : BM026192.D
Acq On : 30 May 2020 03:47
Operator : MA/SJ
Sample : L2820-13
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : CB640

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
1,3-Oxathiolane	4.19	3.3	ng/ul	190131	1	7.47	1159370	20.0
Benzene, chloro-	4.85	2.7	ng/ul	158854	1	7.47	1159370	20.0
o-Cymene	7.62	8.3	ng/ul	482586	1	7.47	1159370	20.0
Indane	7.85	3.3	ng/ul	189771	1	7.47	1159370	20.0
Benzenamine, 3-me...	8.46	7.1	ng/ul	409117	1	7.47	1159370	20.0
Bicyclo[2.2.1]hep...	8.71	4.0	ng/ul	234725	1	7.47	1159370	20.0
Cyclohexanemethan...	9.62	2.5	ng/ul	190000	2	10.23	1508850	20.0
Bicyclo[2.2.1]hep...	9.66	7.6	ng/ul	576580	2	10.23	1508850	20.0
Bicyclo[2.2.1]hep...	9.79	4.1	ng/ul	308397	2	10.23	1508850	20.0
Terpinen-4-ol	10.12	4.1	ng/ul	307922	2	10.23	1508850	20.0
unknown-01	11.08	2.8	ng/ul	209835	2	10.23	1508850	20.0
Phenol, p-tert-bu...	11.60	4.9	ng/ul	371324	2	10.23	1508850	20.0
Benzenamine, 3-ch...	11.75	2.3	ng/ul	171042	2	10.23	1508850	20.0
Naphthalene, 1-me...	12.13	4.6	ng/ul	348461	2	10.23	1508850	20.0
1H-Indole, 3-methyl-	12.97	6.5	ng/ul	1010690	3	14.12	3087520	20.0
1H-Indole, 2,3-di...	13.95	5.6	ng/ul	860011	3	14.12	3087520	20.0
1,1'-Biphenyl, 2-...	14.25	3.3	ng/ul	506026	3	14.12	3087520	20.0
Diethyltoluamide	14.89	14.9	ng/ul	2293640	3	14.12	3087520	20.0
1,1'-Biphenyl, 2,...	15.39	3.7	ng/ul	566686	3	14.12	3087520	20.0
Benzenesulfonamid...	15.65	17.6	ng/ul	1821450	4	16.87	2074000	20.0
2(3H)-Benzothiazo...	15.82	8.9	ng/ul	919380	4	16.87	2074000	20.0
Benzenesulfonamid...	16.16	9.7	ng/ul	1005680	4	16.87	2074000	20.0
Dibenzothiophene	16.69	2.5	ng/ul	256843	4	16.87	2074000	20.0
4H-Cyclopenta[def...	17.94	2.6	ng/ul	268044	4	16.87	2074000	20.0
(DEL) Alkane: Str...	20.82	6.1	ng/ul	711784	5	21.07	2320770	20.0