LSC Area Percent Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM062320\
Data File : BM026519.D
Acq On : 24 Jun 2020 03:30
Operator : JU/CG
Sample : L3069-12
Misc :
ALS Vial : 29 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : CB7H2

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-BM061220MA.M
Title : SVOA CALIBRATION

Signal : TIC: BM026522.D

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.663	127	132	140	rVB	138402	196162	2.55%	0.229%
2	4.087	199	204	214	rVB	84467	134463	1.75%	0.157%
3	6.375	589	593	599	rVB2	49689	71922	0.93%	0.084%
4	6.587	623	629	640	rVB2	163597	344398	4.47%	0.403%
5	6.734	648	654	660	rBV	505069	769285	9.99%	0.900%
6	6.798	660	665	673	rVB5	40620	86905	1.13%	0.102%
7	6.922	680	686	693	rVB	595828	887342	11.52%	1.038%
8	7.046	701	707	714	rVB5	29291	64543	0.84%	0.076%
9	7.381	755	764	774	rBV	534519	875923	11.37%	1.025%
10	7.522	780	788	798	rVB	2201545	3290953	42.72%	3.850%
11	7.740	821	825	831	rVB	142594	214424	2.78%	0.251%
12	8.063	868	880	882	rBV3	52747	132641	1.72%	0.155%
13	8.110	882	888	892	rVV	448666	782091	10.15%	0.915%
14	8.175	892	899	911	rVV	4929185	7703457	100.00%	9.012%
15	8.363	926	931	937	rVB	240553	376644	4.89%	0.441%
16	8.481	945	951	953	rBV2	63386	100942	1.31%	0.118%
17	8.528	953	959	967	rVB	699794	1138079	14.77%	1.331%
18	8.610	967	973	979	rVB	910963	1323724	17.18%	1.549%
19	8.875	1013	1018	1024	rBV5	42900	73715	0.96%	0.086%
20	9.240	1073	1080	1093	rBV2	615031	1187738	15.42%	1.389%
21	9.392	1096	1106	1114	rBV6	97424	323477	4.20%	0.378%
22	9.463	1114	1118	1121	rVV3	50701	81863	1.06%	0.096%
23	9.528	1121	1129	1132	rVV	2564344	4001954	51.95%	4.682%
24	9.563	1132	1135	1143	rVV	1928881	2967943	38.53%	3.472%
25	9.692	1143	1157	1164	rVB3	315677	667926	8.67%	0.781%
26	9.781	1164	1172	1181	rBV	1000769	1685506	21.88%	1.972%
27	9.922	1191	1196	1205	rBV4	62344	167410	2.17%	0.196%
28	10.016	1206	1212	1222	rVV	932233	1495703	19.42%	1.750%
29	10.134	1225	1232	1236	rVV	786391	1333063	17.30%	1.559%
30	10.187	1236	1241	1250	rVV	1299988	2369264	30.76%	2.772%
31	10.287	1250	1258	1272	rVV	773697	1550133	20.12%	1.813%
32	10.404	1273	1278	1282	rVV3	44408	72551	0.94%	0.085%
33	10.498	1289	1294	1300	rBV	132275	213628	2.77%	0.250%
34	10.704	1323	1329	1331	rBV5	35085	66003	0.86%	0.077%

LSC Area Percent Report

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Acq On : 24 Jun 2020 03:30
Operator : JU/CG
Sample : L3069-12
Misc :
ALS Vial : 29 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : CB7H2

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-BM061220MA.M
Title : SVOA CALIBRATION

35	10.739	1331	1335	1342	rVB4	59802	112555	1.46%	0.132%
36	10.992	1371	1378	1391	rVB2	349167	636640	8.26%	0.745%
37	11.510	1457	1466	1473	rBV	163244	302528	3.93%	0.354%
38	11.657	1486	1491	1498	rBV3	162543	289030	3.75%	0.338%
39	11.810	1507	1517	1524	rVB	335459	640770	8.32%	0.750%
40	11.886	1524	1530	1532	rBV2	141016	263803	3.42%	0.309%
41	12.033	1550	1555	1565	rVB2	272682	493130	6.40%	0.577%
42	12.186	1569	1581	1585	rBV5	88639	197310	2.56%	0.231%
43	12.369	1605	1612	1617	rBV8	39524	104752	1.36%	0.123%
44	12.475	1622	1630	1635	rVV5	65009	155790	2.02%	0.182%
45	12.757	1673	1678	1686	rVB4	65336	149010	1.93%	0.174%
46	12.833	1686	1691	1694	rBV2	70761	134477	1.75%	0.157%
47	12.880	1695	1699	1708	rVB	334855	570959	7.41%	0.668%
48	13.080	1730	1733	1742	rVB5	63524	127344	1.65%	0.149%
49	13.186	1746	1751	1752	rBV4	81200	109564	1.42%	0.128%
50	13.351	1775	1779	1785	rVB3	44261	71965	0.93%	0.084%
51	13.404	1785	1788	1793	rBV5	54684	81988	1.06%	0.096%
52	13.469	1793	1799	1803	rBV	1468196	2004730	26.02%	2.345%
53	13.510	1803	1806	1810	rBV	240870	304450	3.95%	0.356%
54	13.592	1816	1820	1823	rBV5	45564	76012	0.99%	0.089%
55	13.675	1830	1834	1837	rBV	248269	294514	3.82%	0.345%
56	13.722	1837	1842	1846	rVV	1655674	2302506	29.89%	2.694%
57	13.763	1846	1849	1861	rVB	272775	505778	6.57%	0.592%
58	13.886	1863	1870	1876	rBV	385601	649725	8.43%	0.760%
59	13.969	1880	1884	1888	rVB2	80348	106282	1.38%	0.124%
60	14.039	1889	1896	1902	rBV2	1751153	2874135	37.31%	3.362%
61	14.098	1902	1906	1912	rVV	852630	1163587	15.10%	1.361%
62	14.174	1916	1919	1925	rVB	98804	136515	1.77%	0.160%
63	14.304	1937	1941	1946	rBV2	114120	217052	2.82%	0.254%
64	14.351	1947	1949	1956	rVV5	74862	140703	1.83%	0.165%
65	14.439	1959	1964	1971	rVB	315415	451097	5.86%	0.528%
66	14.616	1991	1994	1997	rVB	68946	73591	0.96%	0.086%
67	14.816	2021	2028	2042	rVB	1568297	2236206	29.03%	2.616%
68	15.039	2061	2066	2071	rBV	1977663	2764980	35.89%	3.235%
69	15.098	2071	2076	2081	rVB	525723	726740	9.43%	0.850%
70	15.186	2086	2091	2095	rBV	531080	751276	9.75%	0.879%
71	15.316	2109	2113	2117	rBV3	142336	188949	2.45%	0.221%

LSC Area Percent Report

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Acq On : 24 Jun 2020 03:30
Operator : JU/CG
Sample : L3069-12
Misc :
ALS Vial : 29 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : CB7H2

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-BM061220MA.M
Title : SVOA CALIBRATION

72	15.580	2153	2158	2164	rVB	1034389	1376866	17.87%	1.611%
73	15.763	2182	2189	2195	rBV	368958	667968	8.67%	0.781%
74	16.057	2235	2239	2241	rBV2	101701	138516	1.80%	0.162%
75	16.092	2241	2245	2251	rVV	621158	856127	11.11%	1.002%
76	16.157	2253	2256	2263	rVB2	108096	182886	2.37%	0.214%
77	16.374	2289	2293	2295	rBV	76963	102991	1.34%	0.120%
78	16.404	2295	2298	2307	rVB5	158993	237255	3.08%	0.278%
79	16.792	2356	2364	2368	rBV	1565143	2212239	28.72%	2.588%
80	16.833	2368	2371	2376	rVB	917438	1129278	14.66%	1.321%
81	16.892	2376	2381	2386	rBV	2232765	3043260	39.51%	3.560%
82	17.004	2394	2400	2405	rVB	1974878	2533539	32.89%	2.964%
83	17.204	2430	2434	2440	rVB	1038483	1378042	17.89%	1.612%
84	17.310	2449	2452	2456	rBV	55583	73706	0.96%	0.086%
85	17.721	2518	2522	2526	rBV2	131559	205621	2.67%	0.241%
86	17.804	2534	2536	2539	rVB	74345	72152	0.94%	0.084%
87	17.862	2543	2546	2555	rVB3	85756	144135	1.87%	0.169%
88	18.862	2712	2716	2720	rVB	213858	257587	3.34%	0.301%
89	18.957	2730	2732	2738	rVB	124711	146048	1.90%	0.171%
90	19.080	2750	2753	2757	rBV3	86414	114018	1.48%	0.133%
91	19.198	2768	2773	2781	rBV	2490831	3503961	45.49%	4.099%
92	19.274	2781	2786	2796	rVB	363715	546259	7.09%	0.639%
93	19.615	2842	2844	2849	rVB	69448	75419	0.98%	0.088%
94	19.721	2857	2862	2866	rBV	198965	323339	4.20%	0.378%
95	20.751	3033	3037	3045	rVB2	755450	889978	11.55%	1.041%
96	21.003	3076	3080	3089	rVB	1943418	2395452	31.10%	2.802%
97	22.962	3408	3413	3419	rVV	859882	1363073	17.69%	1.595%
98	23.021	3420	3423	3428	rVB	173626	242936	3.15%	0.284%
99	23.092	3430	3435	3443	rVB	1142466	1852860	24.05%	2.168%
100	23.603	3519	3522	3529	rVB	159523	253308	3.29%	0.296%

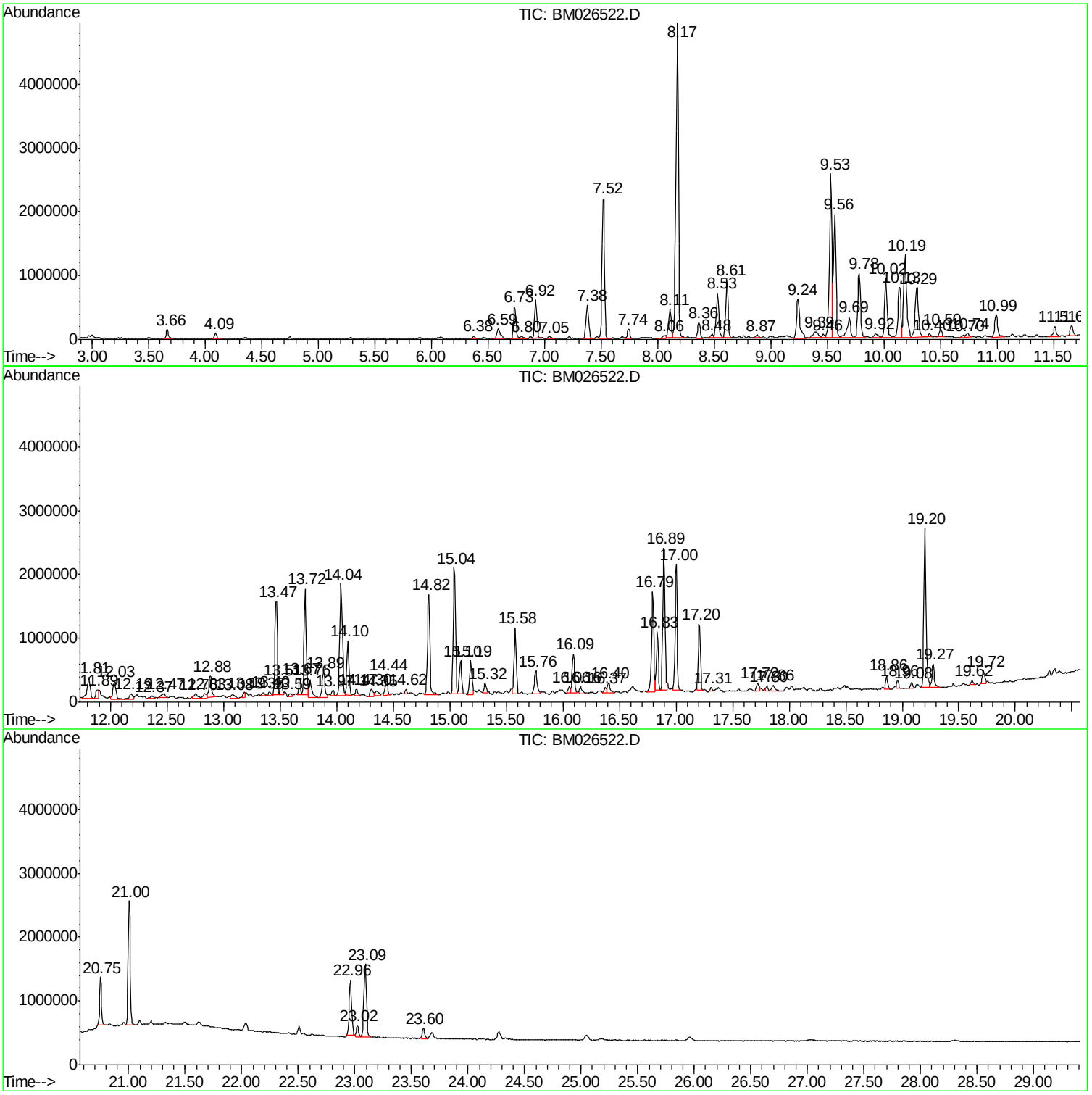
Sum of corrected areas: 85481037

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM062320\
Data File : BM026519.D
Acq On : 24 Jun 2020 03:30
Operator : JU/CG
Sample : L3069-12
Misc :
ALS Vial : 29 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : CB7H2

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM061220MA.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\BM062320\
Data File : BM026519.D
Acq On : 24 Jun 2020 03:30
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Instrument : BNA_M
ClientSampleId : CB7H2

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM061220MA.M
Quant Title : SVOA CALIBRATION

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

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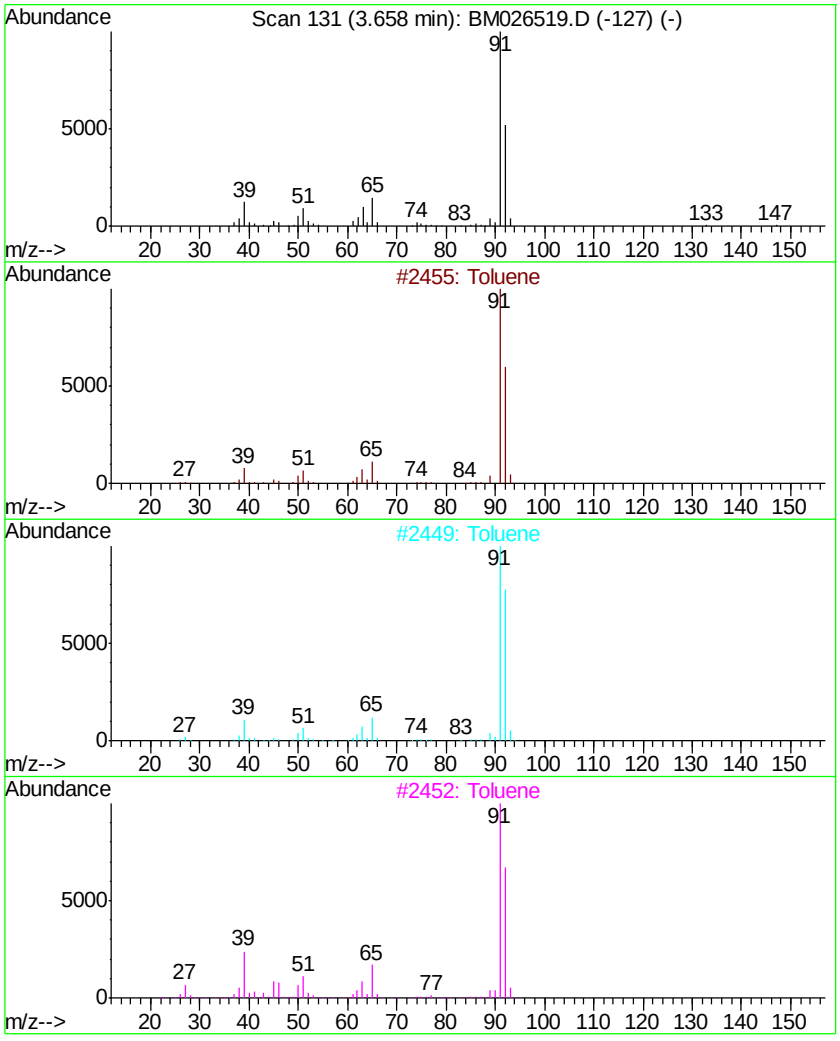
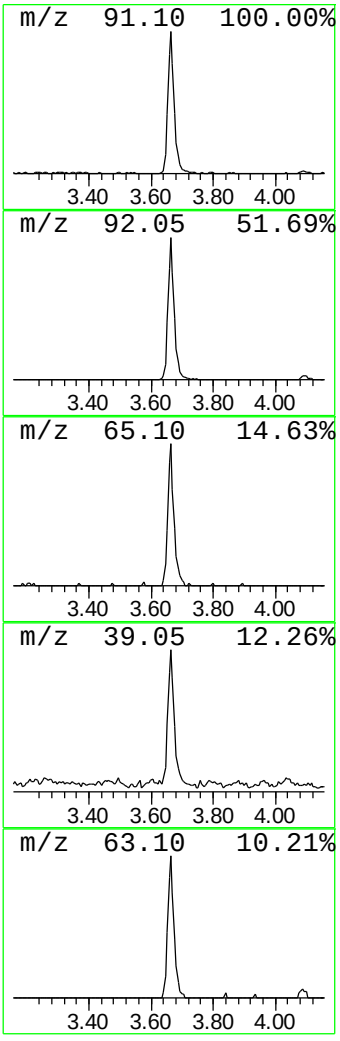
Peak Number 1 Toluene Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.66	4.48 ng/ul	196162	1,4-Dichlorobenzene-d4	7.38

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Toluene	92	C7H8	000108-88-3	94
2	Toluene	92	C7H8	000108-88-3	91
3	Toluene	92	C7H8	000108-88-3	90
4	Spiro[2,4]hepta-4,6-diene	92	C7H8	000765-46-8	90
5	1,3,5-Cycloheptatriene	92	C7H8	000544-25-2	90





Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM062320\
Data File : BM026519.D
Acq On : 24 Jun 2020 03:30
Operator : JU/CG
Sample : L3069-12
Misc :
ALS Vial : 29 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : CB7H2

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM061220MA.M
Quant Title : SVOA CALIBRATION

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

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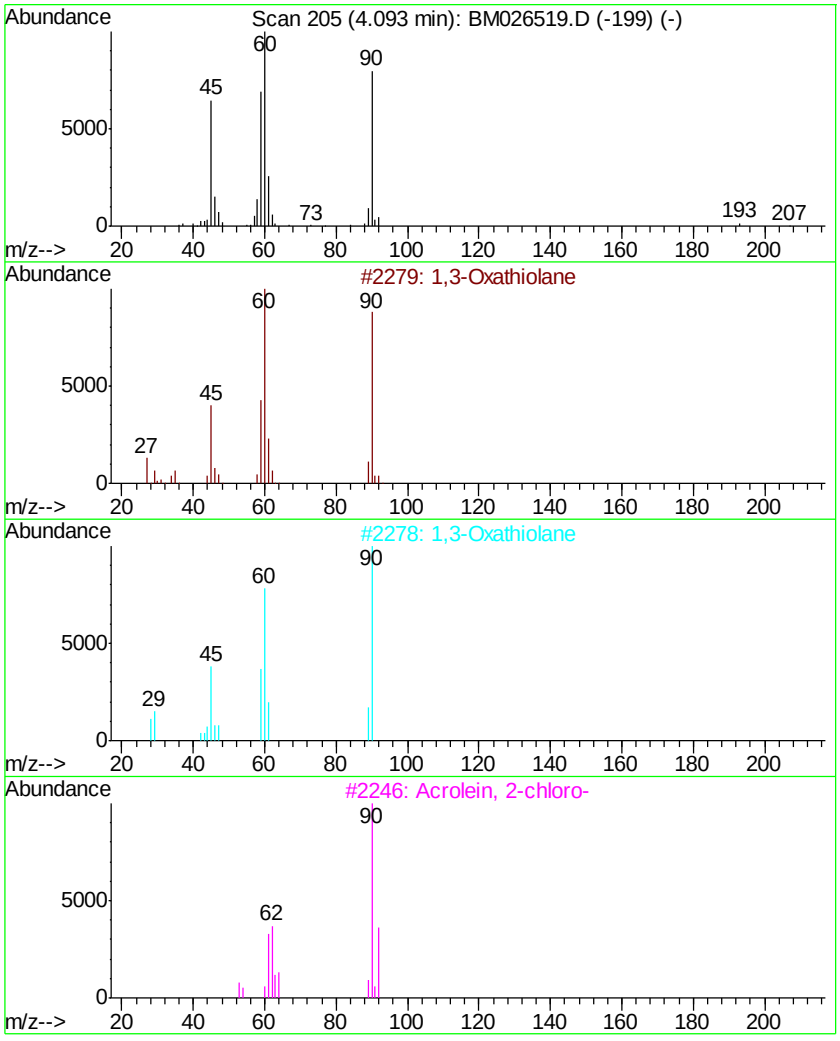
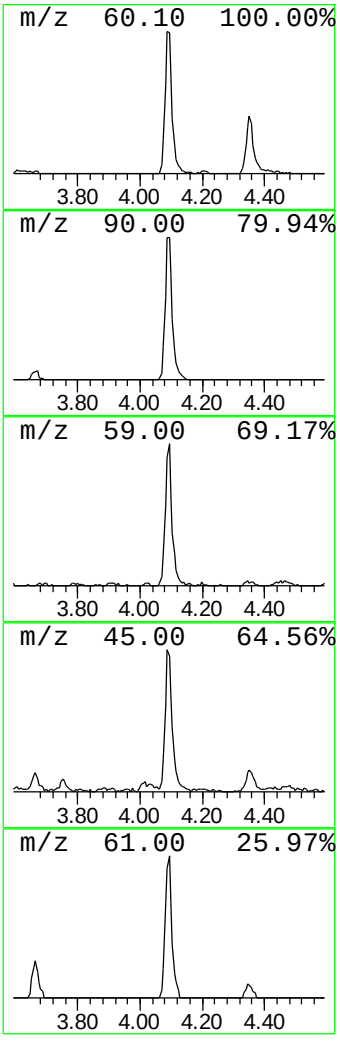
Peak Number 2 1,3-Oxathiolane Concentration Rank 25

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.09	3.07 ng/ul	134463	1,4-Dichlorobenzene-d4	7.38

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,3-Oxathiolane	90	C3H6OS	002094-97-5	91
2	1,3-Oxathiolane	90	C3H6OS	002094-97-5	64
3	Acrolein, 2-chloro-	90	C3H3ClO	000683-51-2	59
4	3-Thietanol	90	C3H6OS	010304-16-2	53
5	Noxytiolin	120	C3H8N2OS	015599-39-0	42





Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM062320\
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Sample : L3069-12
Misc :
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Instrument : BNA_M
ClientSampleId : CB7H2

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM061220MA.M
Quant Title : SVOA CALIBRATION

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

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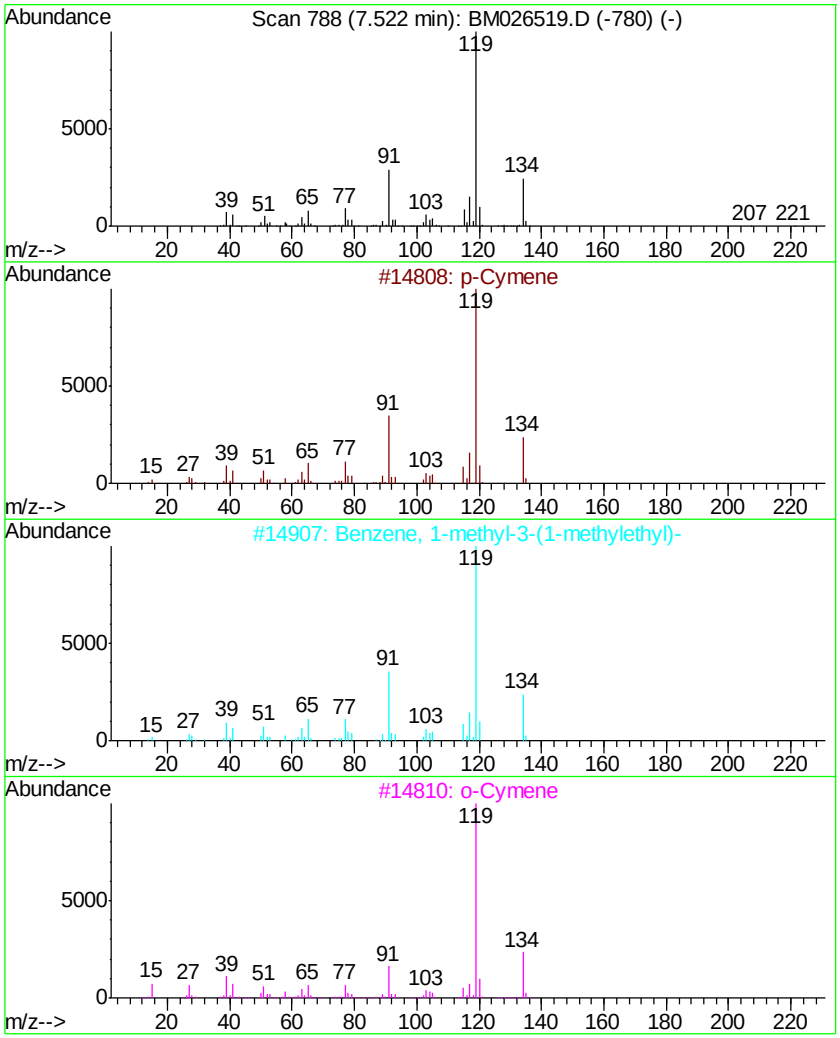
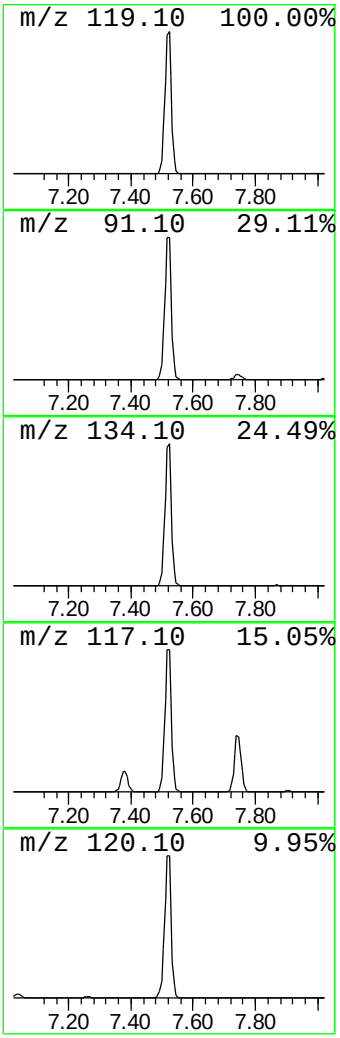
Peak Number 3 p-Cymene Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.52	75.14 ng/ul	3290950	1,4-Dichlorobenzene-d4	7.38

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	o-Cymene	134	C10H14	000527-84-4	94
4	o-Cymene	134	C10H14	000527-84-4	94
5	Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	91





Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM062320\
Data File : BM026519.D
Acq On : 24 Jun 2020 03:30
Operator : JU/CG
Sample : L3069-12
Misc :
ALS Vial : 29 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : CB7H2

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM061220MA.M
Quant Title : SVOA CALIBRATION

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

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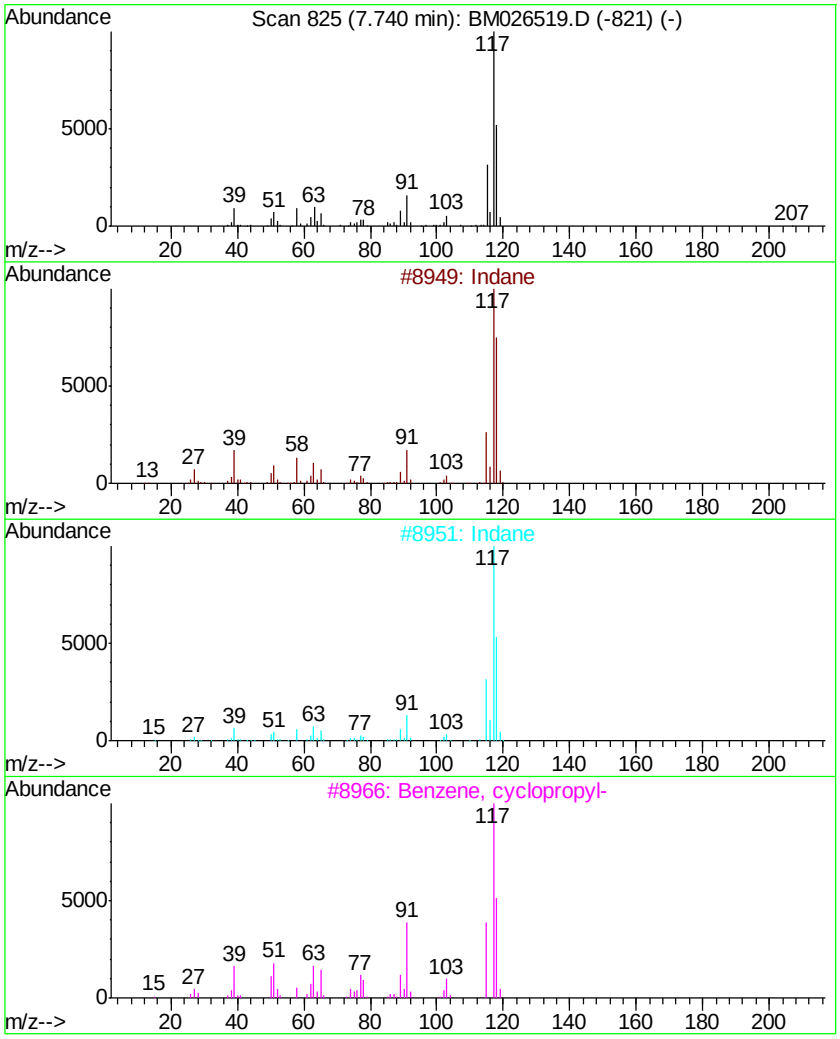
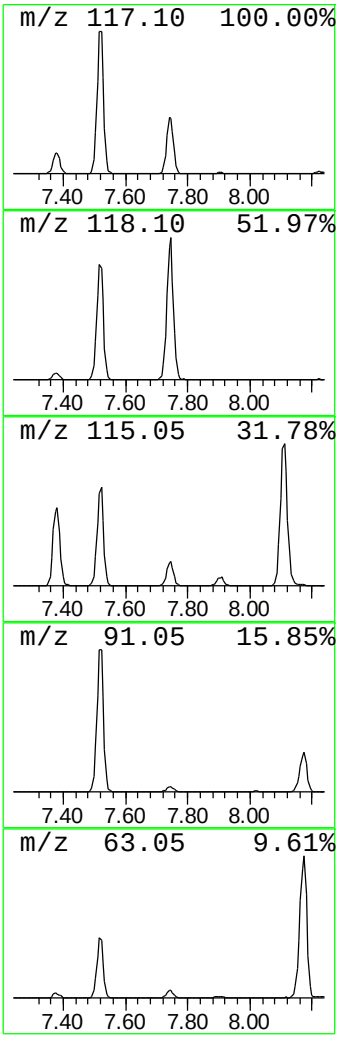
Peak Number 4 Indane Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.74	4.90 ng/ul	214424	1,4-Dichlorobenzene-d4	7.38

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indane	118	C9H10	000496-11-7	96
2	Indane	118	C9H10	000496-11-7	94
3	Benzene, cyclopropyl-	118	C9H10	000873-49-4	93
4	Benzene, 1-ethenyl-2-methyl-	118	C9H10	000611-15-4	90
5	Tetracyclo[3.3.1.0(2,8).0(4,6)]-...	118	C9H10	1000191-13-7	74





Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM062320\
Data File : BM026519.D
Acq On : 24 Jun 2020 03:30
Operator : JU/CG
Sample : L3069-12
Misc :
ALS Vial : 29 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : CB7H2

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM061220MA.M
Quant Title : SVOA CALIBRATION

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

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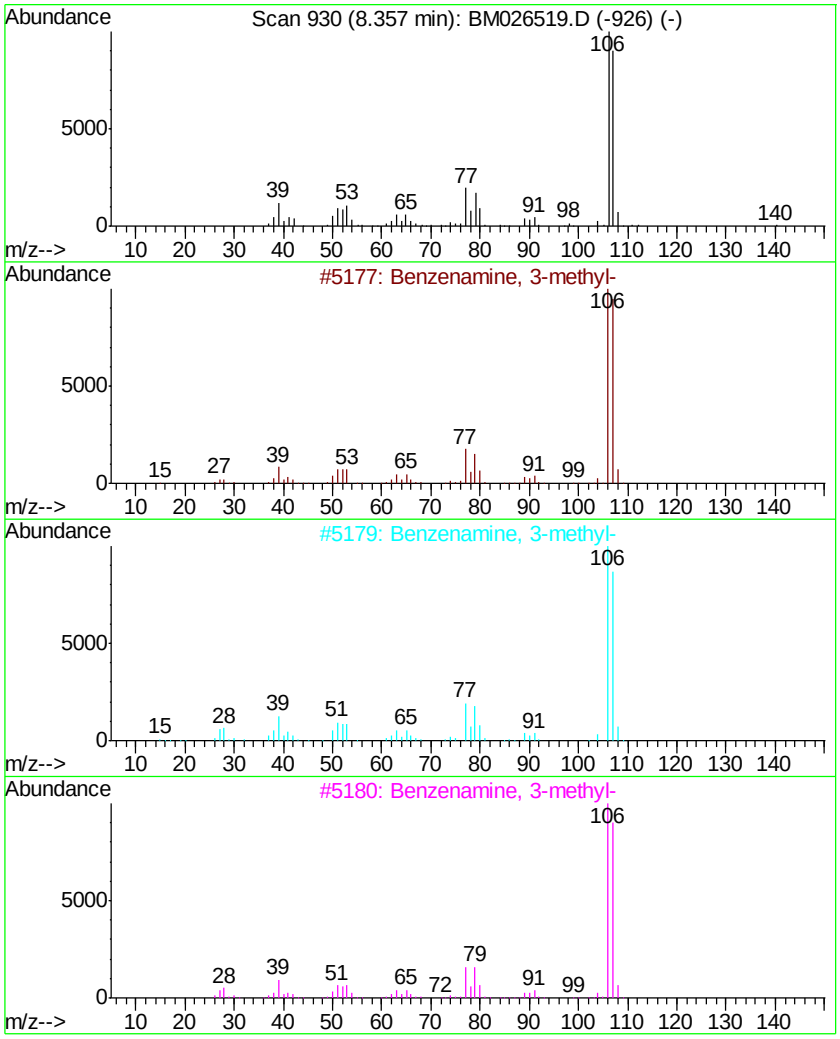
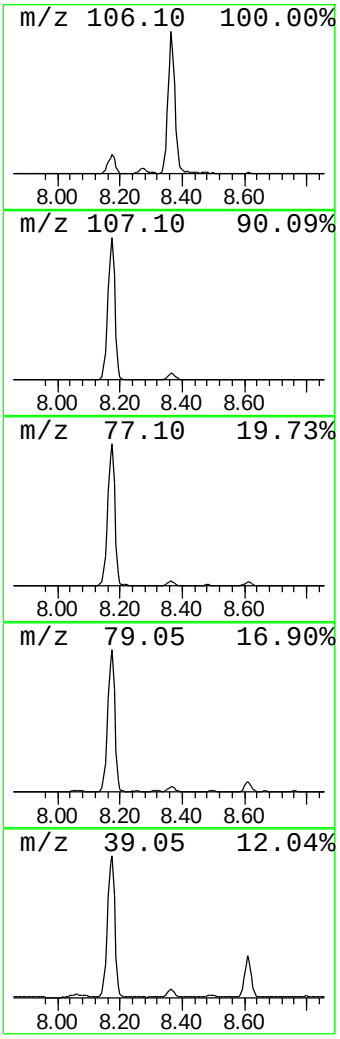
Peak Number 5 Benzenamine, 3-methyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.36	8.60 ng/ul	376644	1,4-Dichlorobenzene-d4	7.38

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	96
3	Benzenamine, 3-methyl-	107	C7H9N	000108-44-1	94
4	Benzenamine, 3-methyl-	107	C7H9N	000108-44-1	94
5	p-Aminotoluene	107	C7H9N	000106-49-0	91





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Acq On : 24 Jun 2020 03:30
Operator : JU/CG
Sample : L3069-12
Misc :
ALS Vial : 29 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : CB7H2

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Quant Title : SVOA CALIBRATION

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

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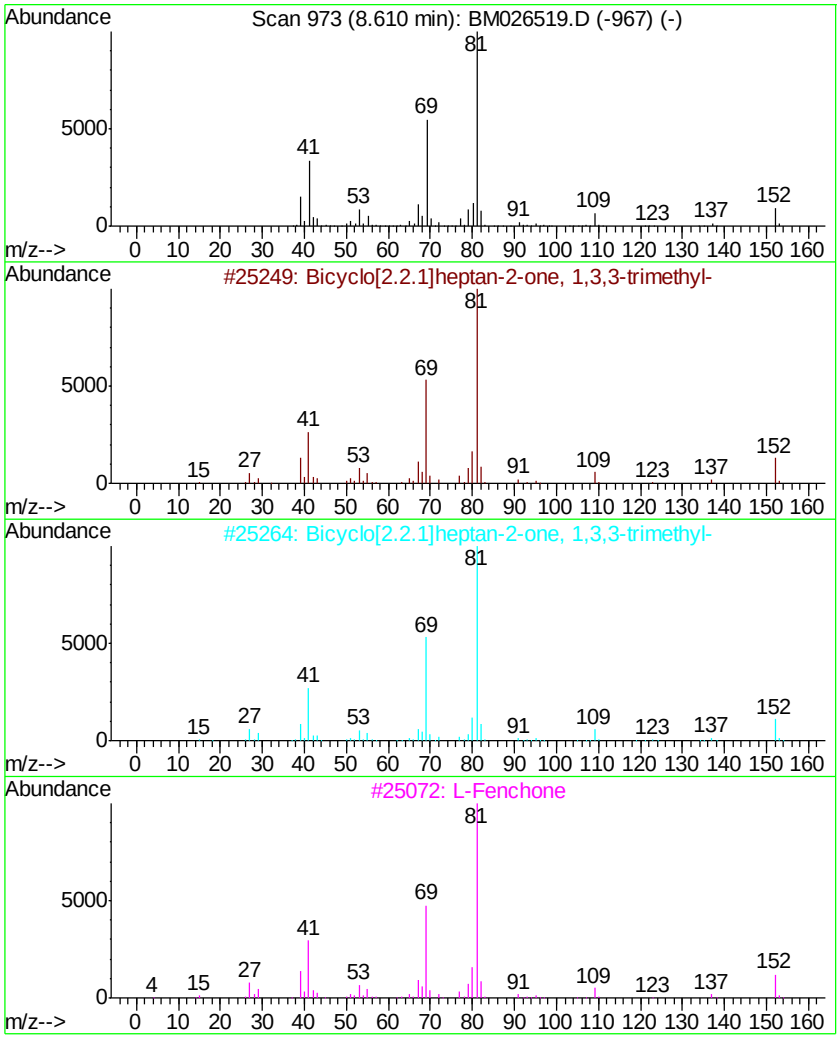
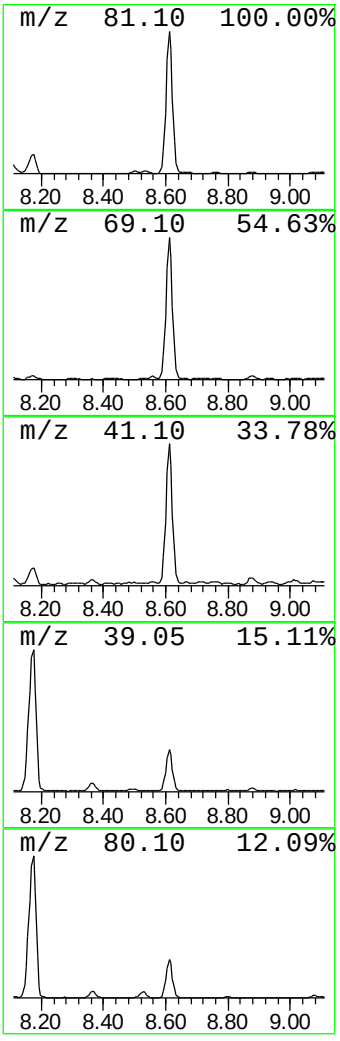
Peak Number 6 Bicyclo[2.2.1]heptan-2-one,... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.61	30.22 ng/ul	1323720	1,4-Dichlorobenzene-d4	7.38

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	D-Fenchone	152	C10H16O	004695-62-9	91
5	L-Fenchone	152	C10H16O	007787-20-4	91





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ClientSampleId : CB7H2

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Quant Title : SVOA CALIBRATION

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

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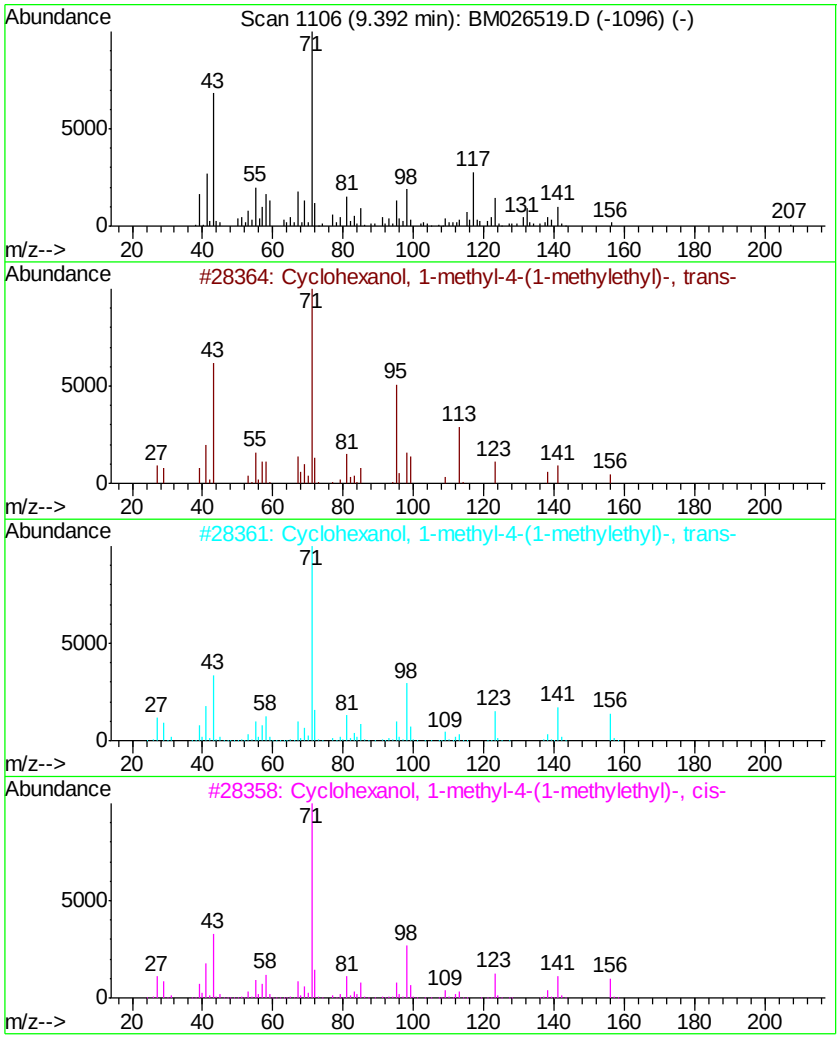
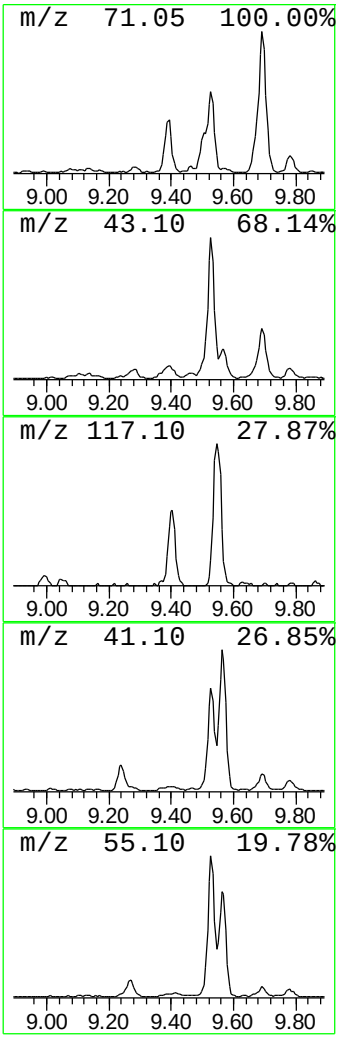
Peak Number 7 Cyclohexanol, 1-methyl-4-(1... Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.39	4.85 ng/ul	323477	Naphthalene-d8	10.14

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclohexanol, 1-methyl-4-(1-meth...	156	C10H20O	003901-93-7	59
2	Cyclohexanol, 1-methyl-4-(1-meth...	156	C10H20O	003901-93-7	53
3	Cyclohexanol, 1-methyl-4-(1-meth...	156	C10H20O	003901-95-9	53
4	1-Methylcyclooctanol	142	C9H18O	059123-41-0	47
5	2-Nonen-1-ol, 2-methyl-	156	C10H20O	091008-40-1	47





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Instrument : BNA_M
ClientSampleId : CB7H2

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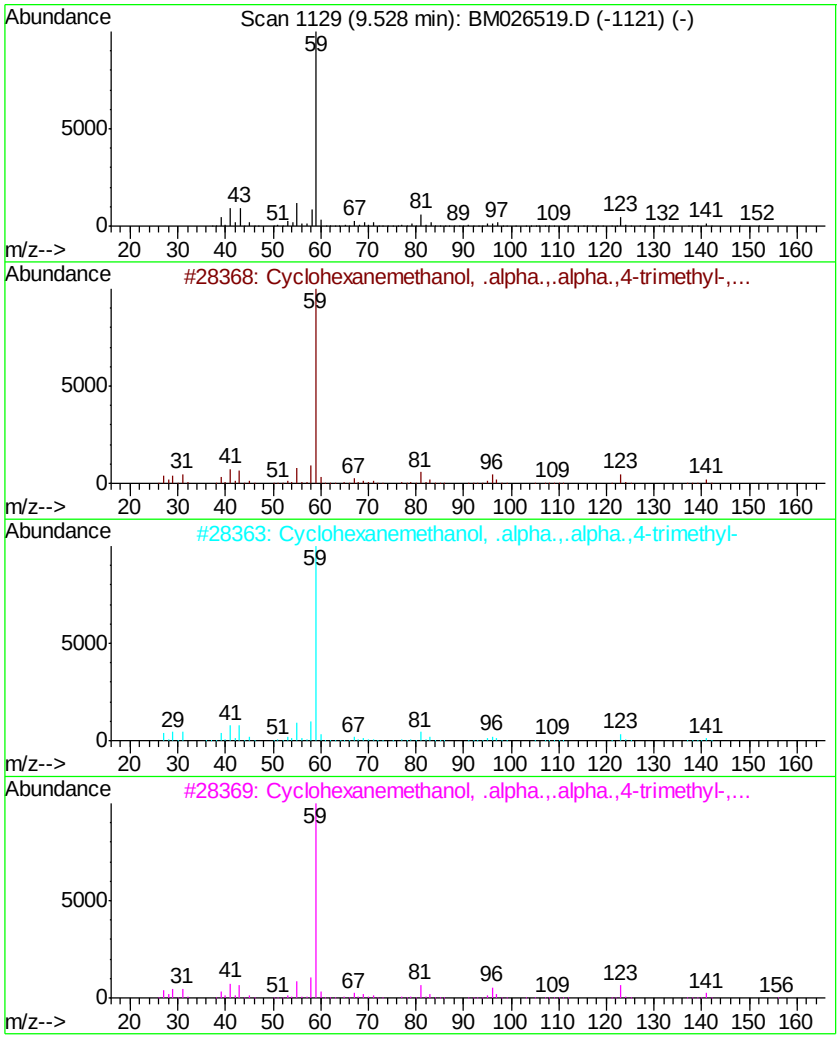
Quant Title : SVOA CALIBRATION

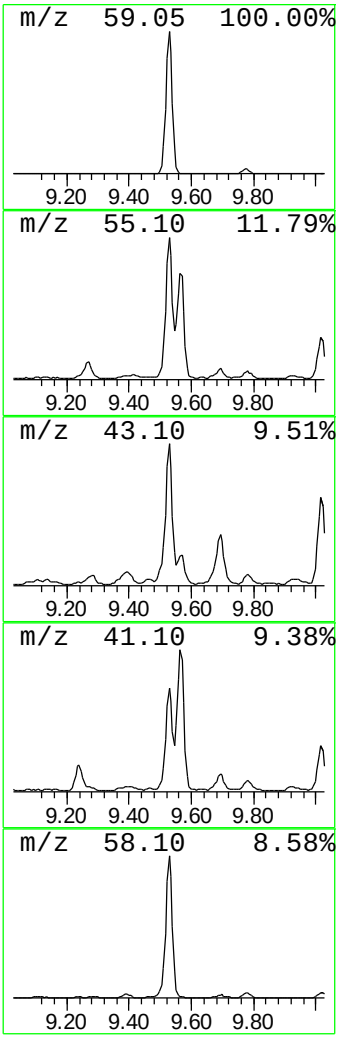
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 8 Cyclohexanemethanol, .alpha... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.53	60.04 ng/ul	4001950	Naphthalene-d8	10.14

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclohexanemethanol, .alpha.,.al...	156	C10H20O	007322-63-6	83
2	Cyclohexanemethanol, .alpha.,.al...	156	C10H20O	000498-81-7	78
3	Cyclohexanemethanol, .alpha.,.al...	156	C10H20O	005114-00-1	74
4	Cyclohexanemethanol, .alpha.,.al...	156	C10H20O	000498-81-7	47
5	2-Hydroxy-2,5-dimethyl-hept-6-en...	156	C9H16O2	1000192-55-8	45





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ClientSampleId : CB7H2

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Quant Title : SVOA CALIBRATION

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

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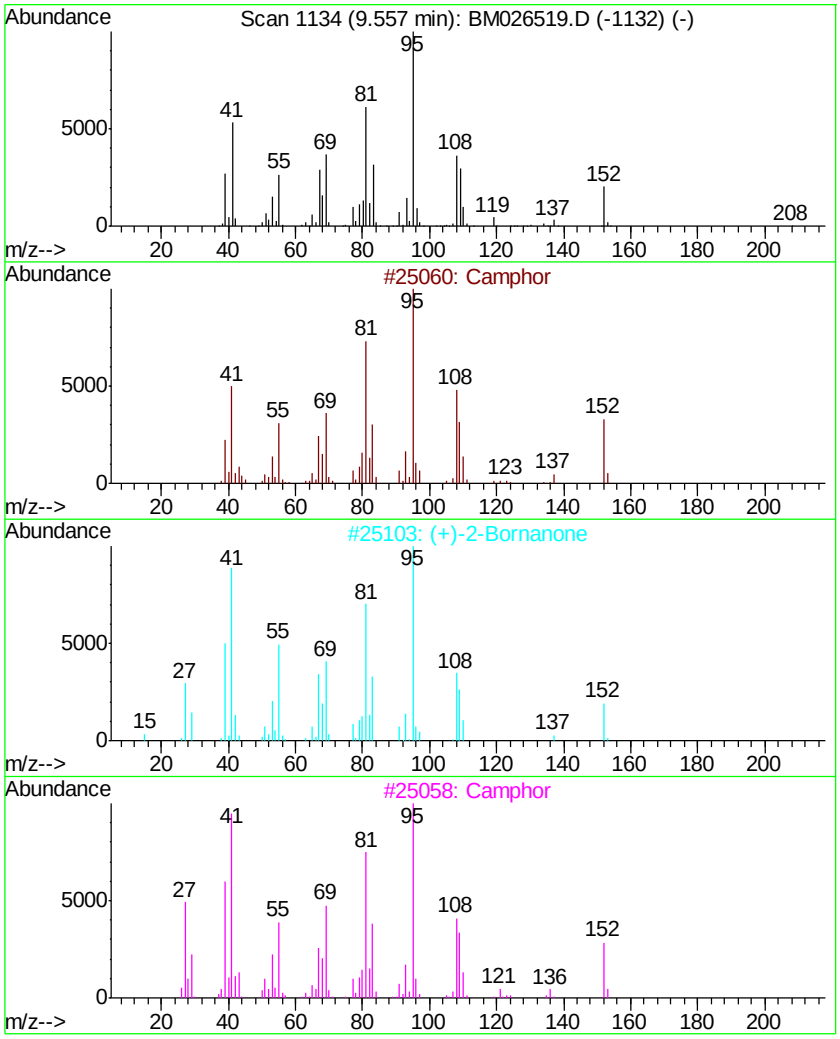
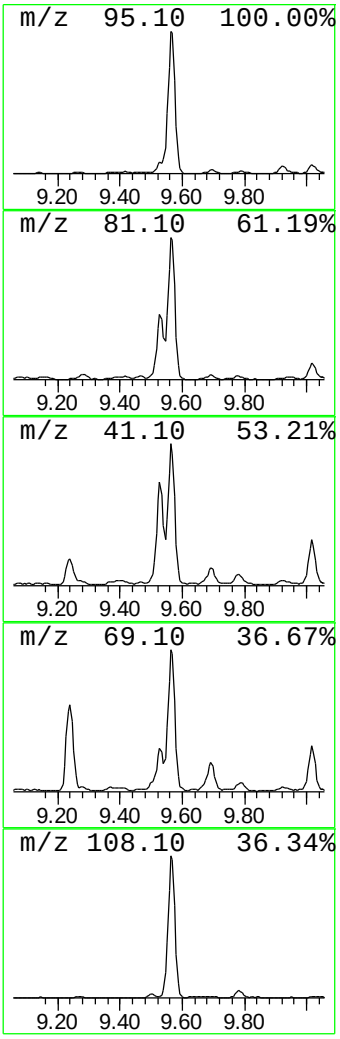
Peak Number 9 Camphor Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.56	44.53 ng/ul	2967940	Naphthalene-d8	10.14

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





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Misc :
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Quant Title : SVOA CALIBRATION

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

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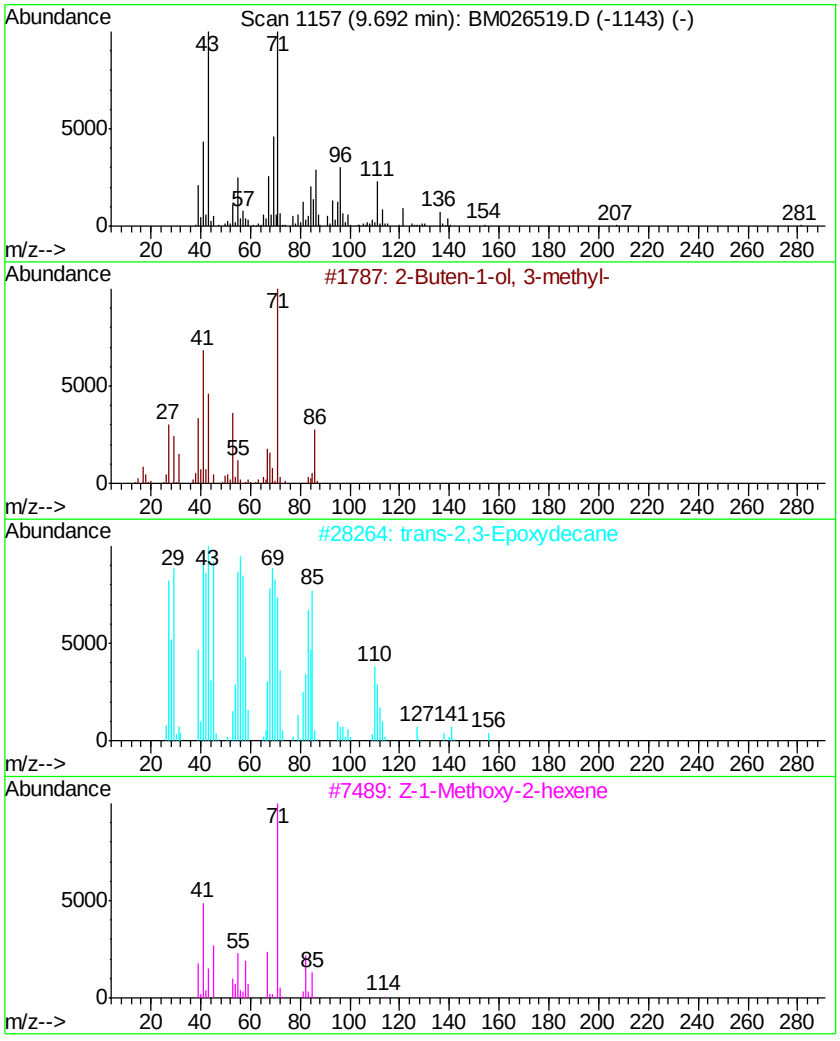
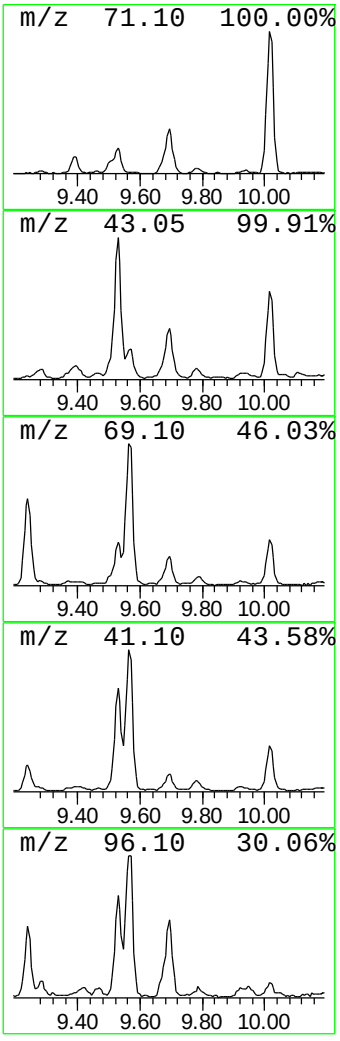
Peak Number 10 unknown-01 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.69	10.02 ng/ul	667926	Naphthalene-d8	10.14

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Buten-1-ol, 3-methyl-	86	C5H10O	000556-82-1	47
2	trans-2,3-Epoxydecane	156	C10H20O	054125-39-2	47
3	Z-1-Methoxy-2-hexene	114	C7H14O	1000279-60-9	45
4	3-Methoxybut-1-ene	86	C5H10O	017351-24-5	43
5	2-Butene, 1-methoxy-, (E)-	86	C5H10O	010034-14-7	38





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ClientSampleId : CB7H2

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Quant Title : SVOA CALIBRATION

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

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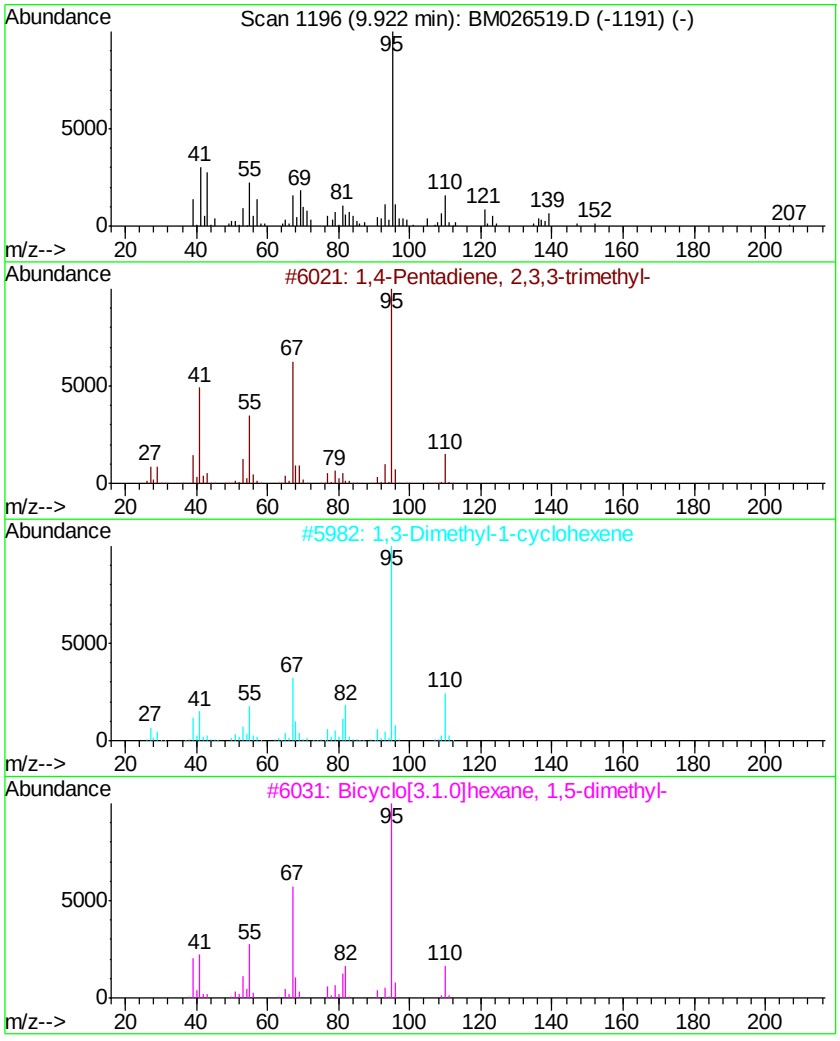
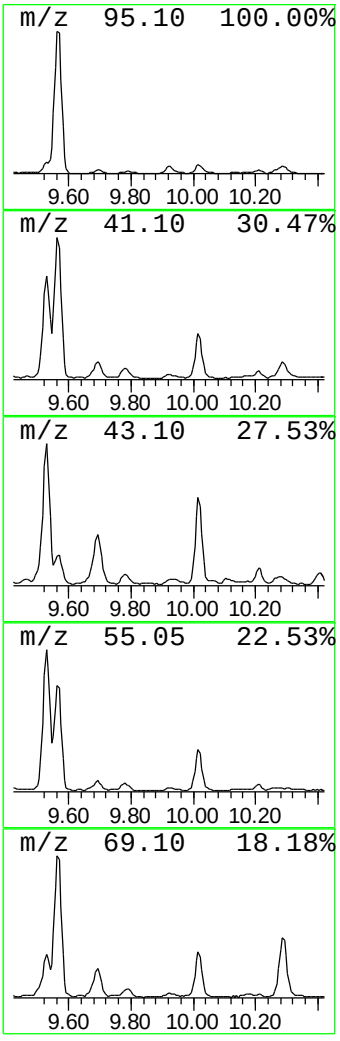
Peak Number 11 1,4-Pentadiene, 2,3,3-trime... Concentration Rank 29

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.92	2.51 ng/ul	167410	Naphthalene-d8	10.14

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,4-Pentadiene, 2,3,3-trimethyl-	110	C8H14	000756-02-5	68
2	1,3-Dimethyl-1-cyclohexene	110	C8H14	002808-76-6	64
3	Bicyclo[3.1.0]hexane, 1,5-dimethyl-	110	C8H14	1000142-17-5	64
4	1-(1-Propynyl)cyclohexanol	138	C9H14O	000697-37-0	59
5	1,5-Hexadiene, 2,5-dimethyl-	110	C8H14	000627-58-7	59





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Instrument : BNA_M
ClientSampleId : CB7H2

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Quant Title : SVOA CALIBRATION

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

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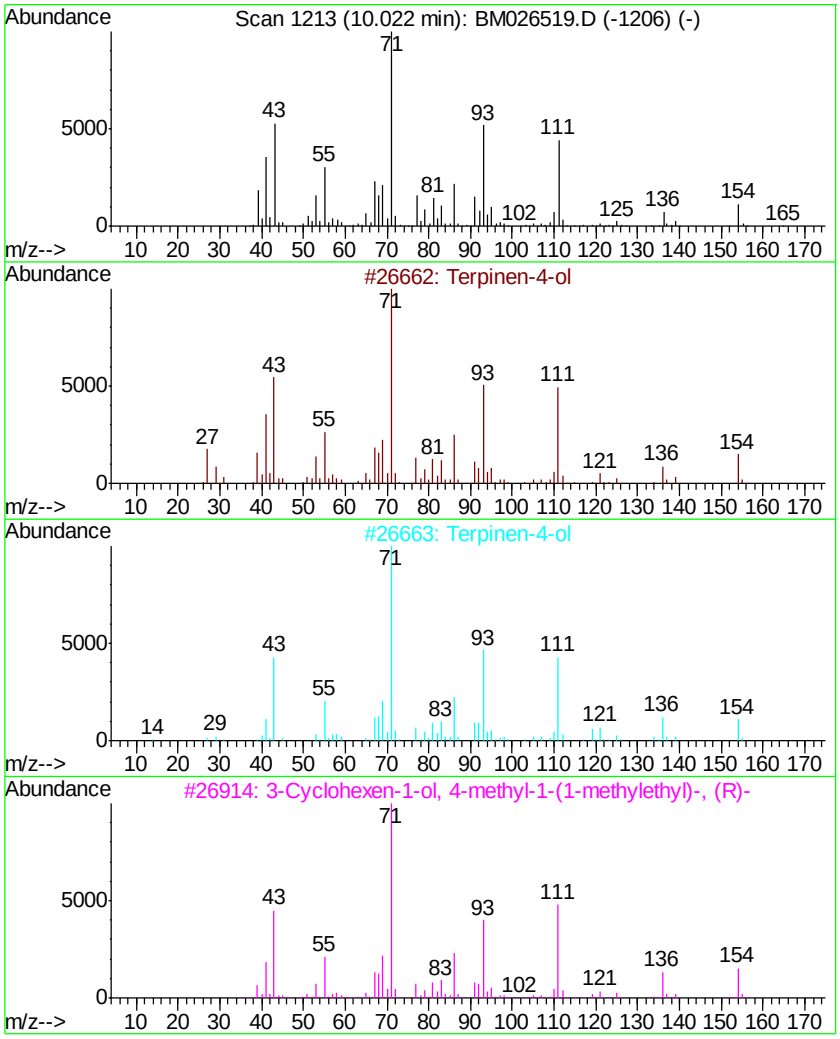
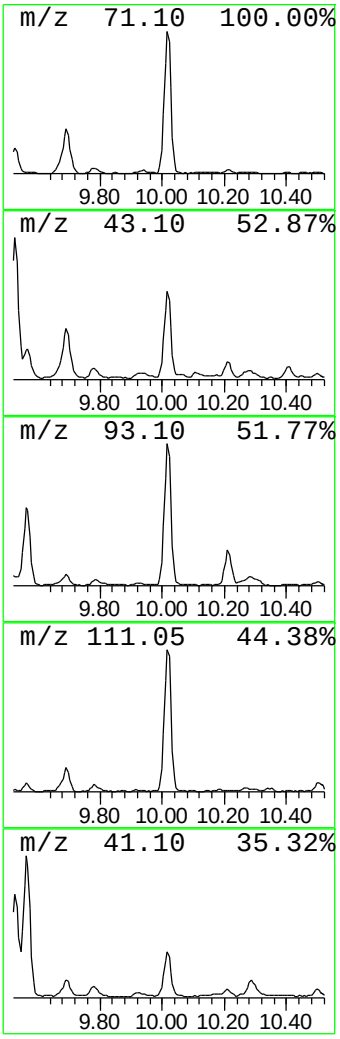
Peak Number 12 Terpinen-4-ol Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.02	22.44 ng/ul	1495700	Naphthalene-d8	10.14

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	91
3	3-Cyclohexen-1-ol, 4-methyl-1-(1...	154	C10H18O	020126-76-5	90
4	Terpinen-4-ol	154	C10H18O	000562-74-3	81
5	3-Cyclohexen-1-ol, 4-methyl-1-(1...	154	C10H18O	020126-76-5	81





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Quant Title : SVOA CALIBRATION

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

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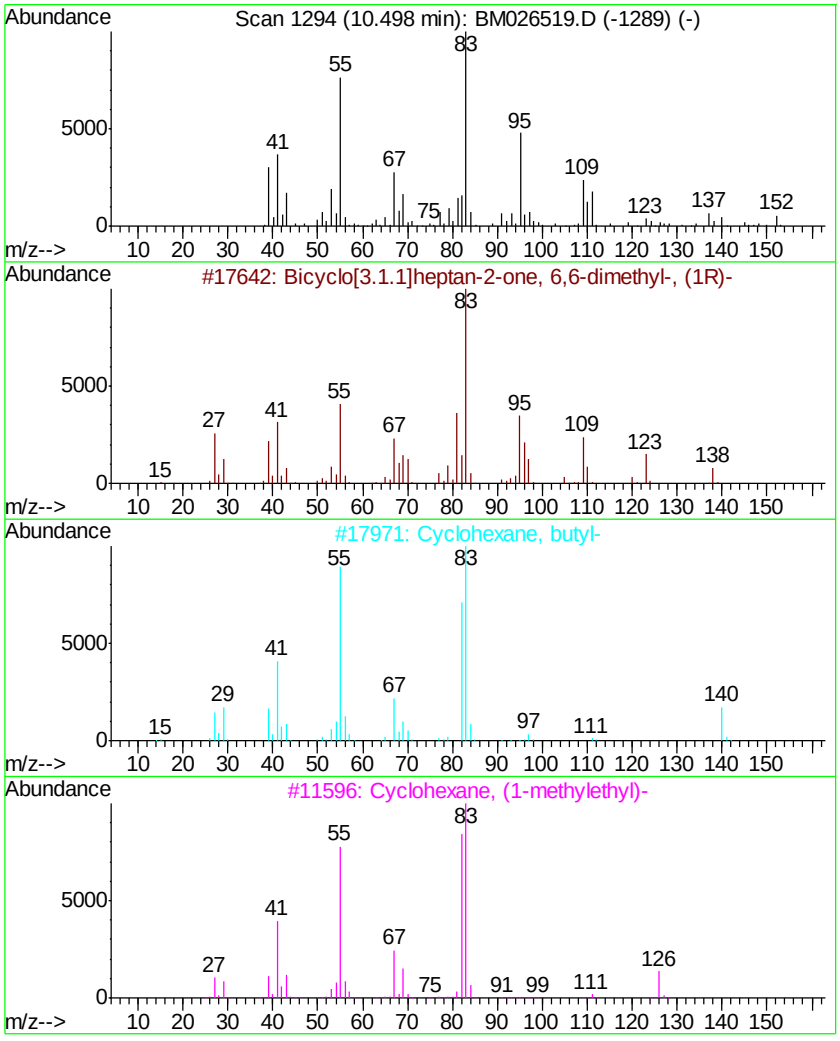
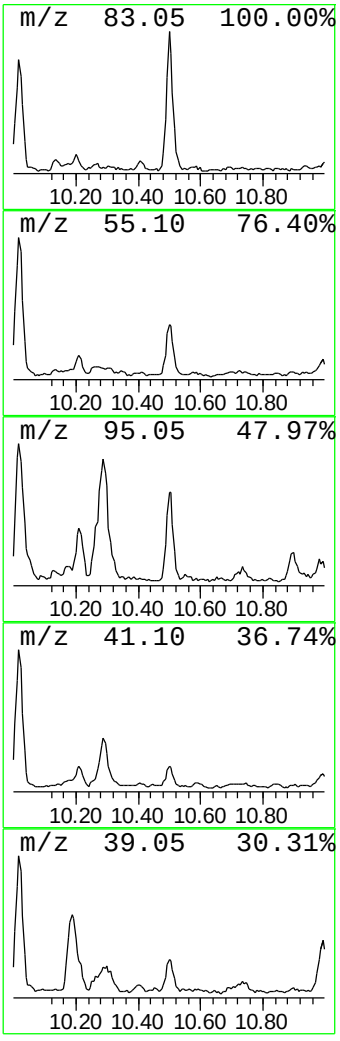
Peak Number 13 Bicyclo[3.1.1]heptan-2-one,... Concentration Rank 24

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.50	3.21 ng/ul	213628	Naphthalene-d8	10.14

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Bicyclo[3.1.1]heptan-2-one, 6,6-...	138	C9H14O	038651-65-9	72
2	Cyclohexane, butyl-	140	C10H20	001678-93-9	43
3	Cyclohexane, (1-methylethyl)-	126	C9H18	000696-29-7	43
4	1-Propanone, 1-cyclohexyl-	140	C9H16O	001123-86-0	43
5	1,6-Octadiene, 5,7-dimethyl-, (R)-	138	C10H18	085006-04-8	40





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Quant Title : SVOA CALIBRATION

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

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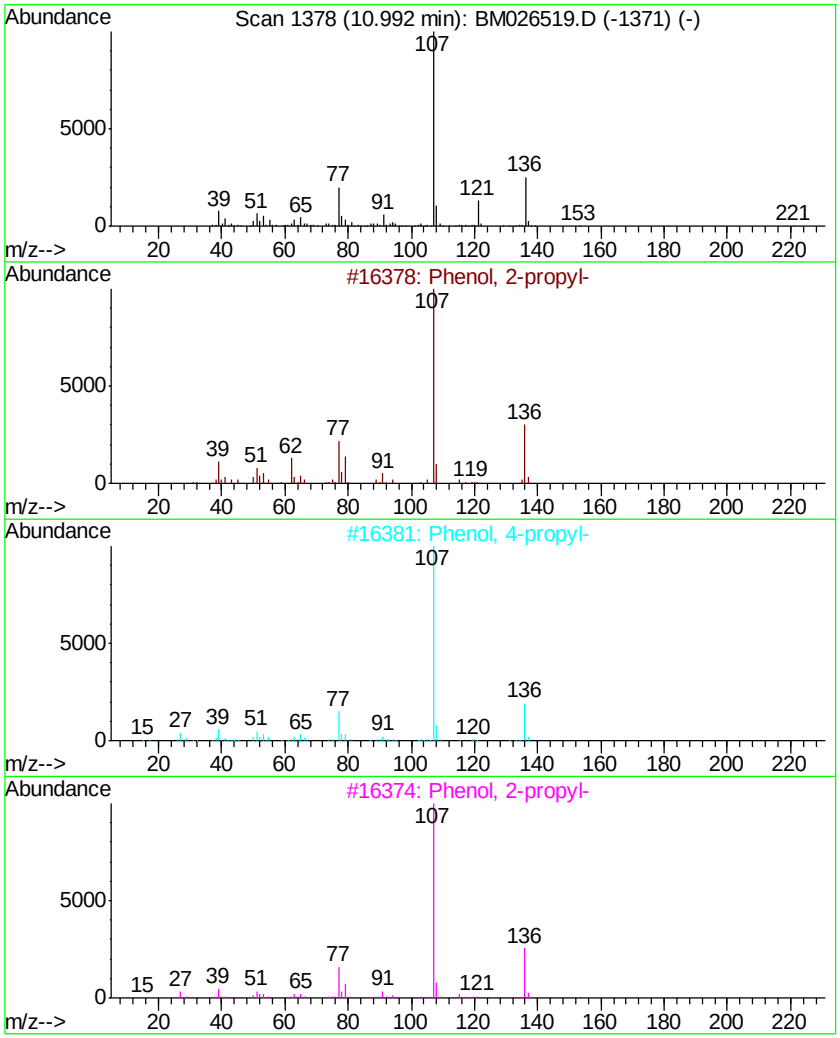
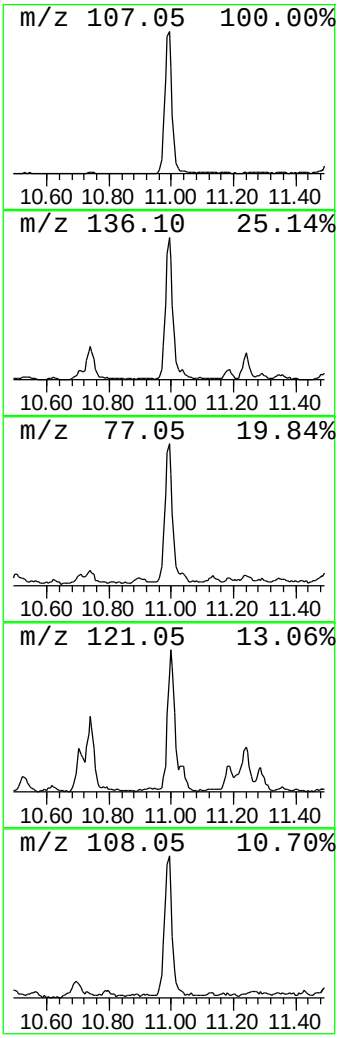
Peak Number 14 Phenol, 2-propyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.99	9.55 ng/ul	636640	Naphthalene-d8	10.14

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, 2-propyl-	136	C9H12O	000644-35-9	87
2	Phenol, 4-propyl-	136	C9H12O	000645-56-7	87
3	Phenol, 2-propyl-	136	C9H12O	000644-35-9	87
4	Cyclohexene, 1-ethyl-6-ethylidene-	136	C10H16	061141-57-9	80
5	Phenol, 2-propyl-	136	C9H12O	000644-35-9	80





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Quant Title : SVOA CALIBRATION

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

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Peak Number 15 Phenol, p-tert-butyl- Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.51	4.54 ng/ul	302528	Naphthalene-d8	10.14

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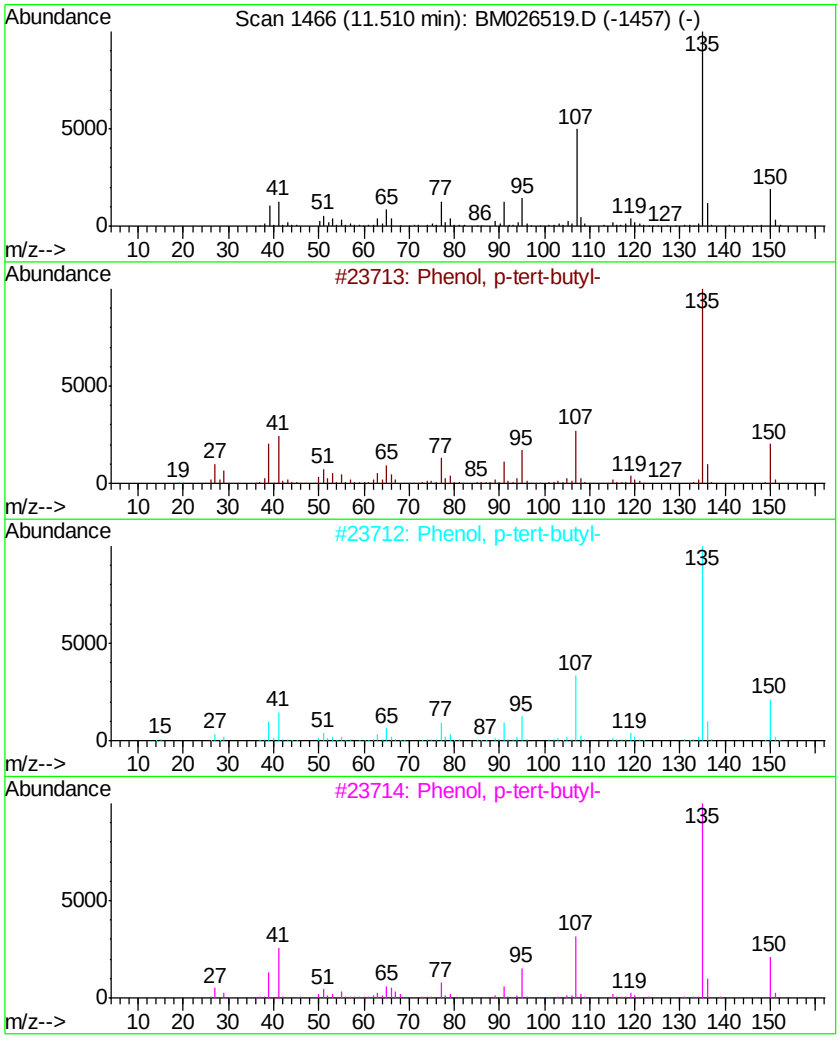

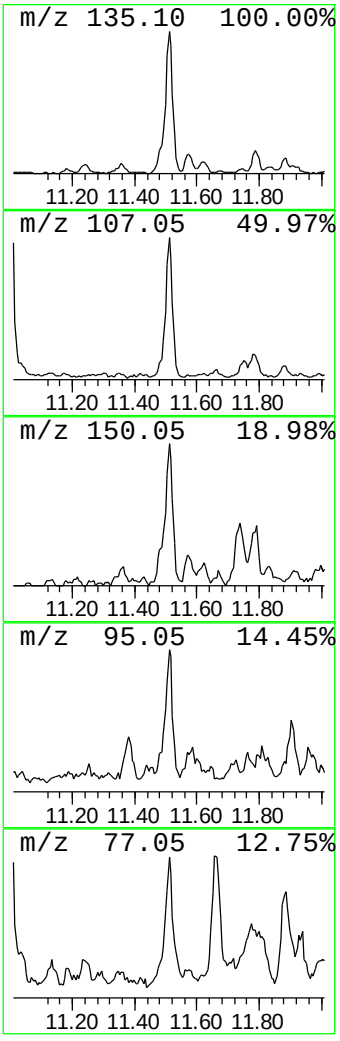

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Quant Title : SVOA CALIBRATION

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

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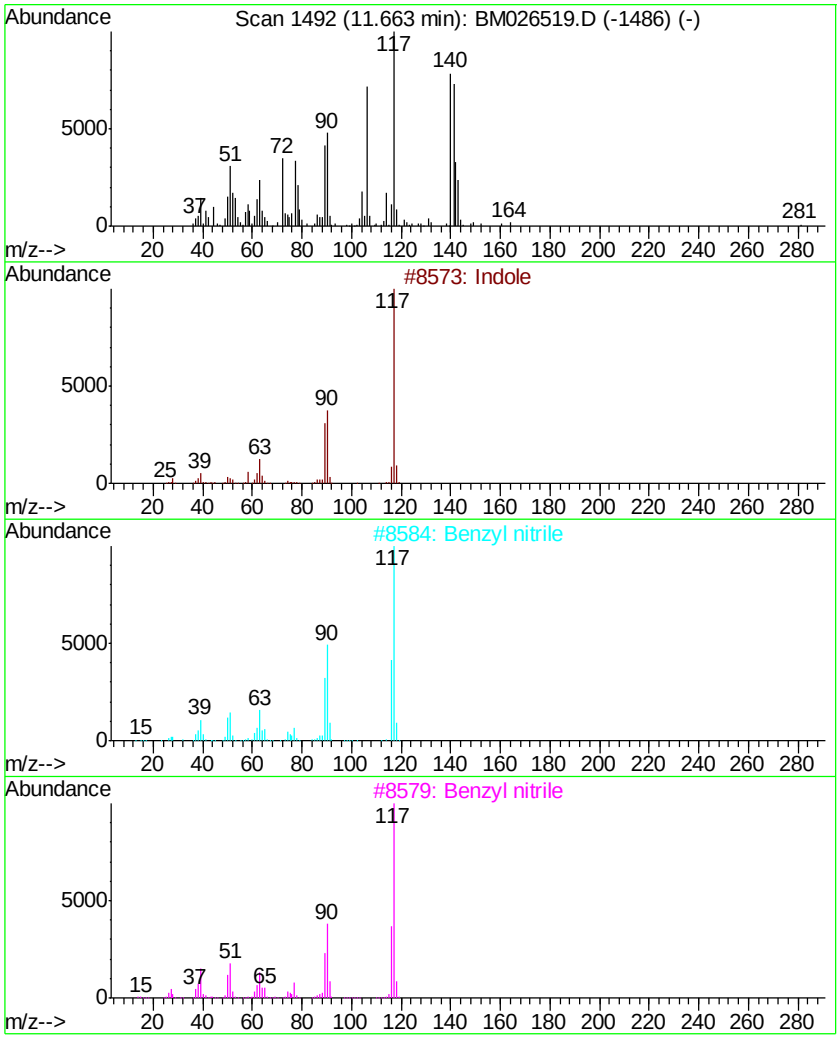
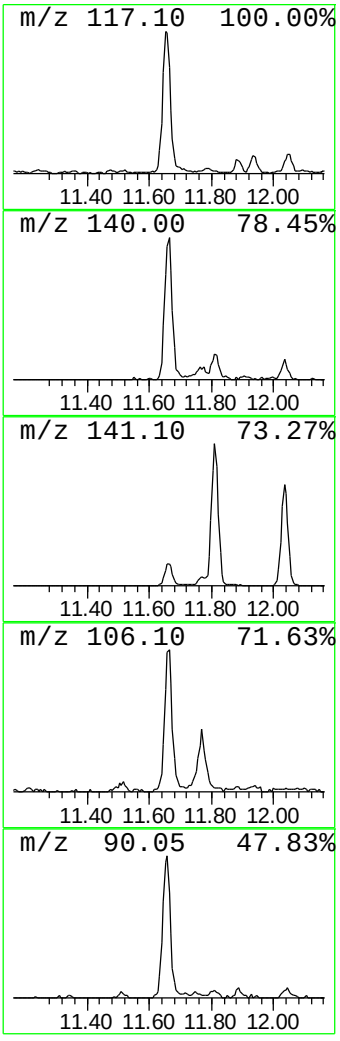
Peak Number 16 Indole Concentration Rank 21

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.66	4.34 ng/ul	289030	Naphthalene-d8	10.14

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indole	117	C8H7N	000120-72-9	70
2	Benzyl nitrile	117	C8H7N	000140-29-4	70
3	Benzyl nitrile	117	C8H7N	000140-29-4	55
4	Benzonitrile, 2-methyl-	117	C8H7N	000529-19-1	55
5	Indole	117	C8H7N	000120-72-9	55





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ClientSampleId : CB7H2

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM061220MA.M
Quant Title : SVOA CALIBRATION

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

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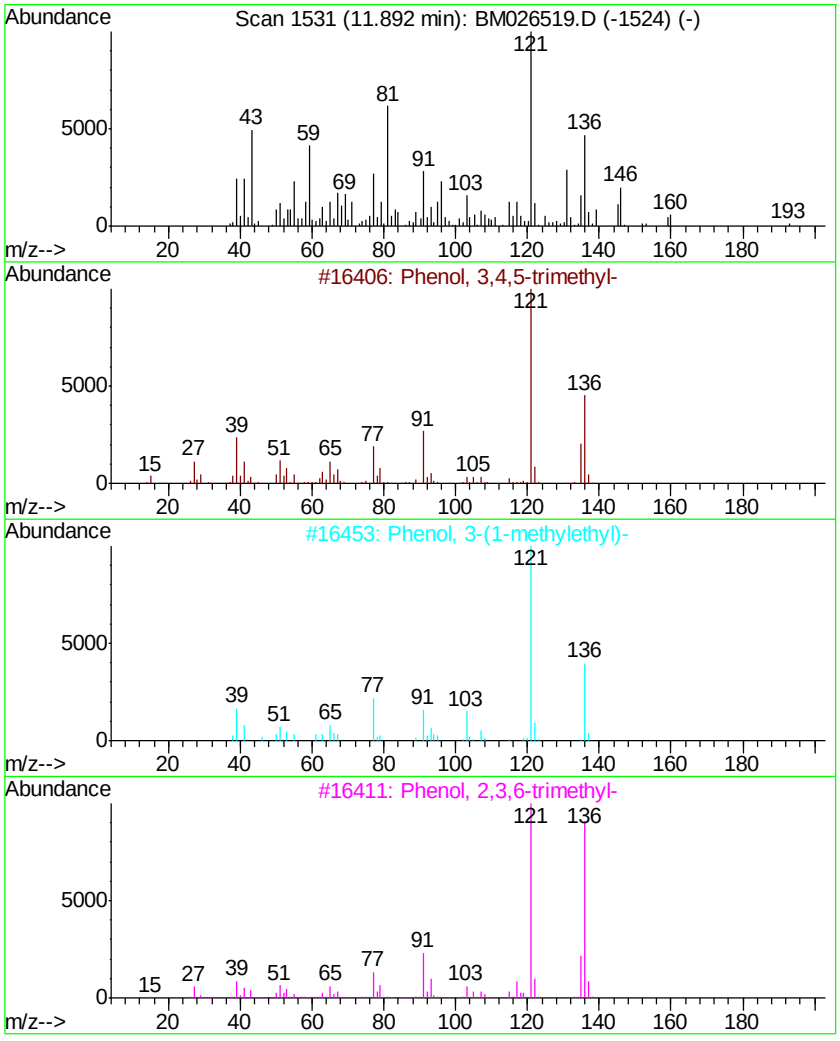
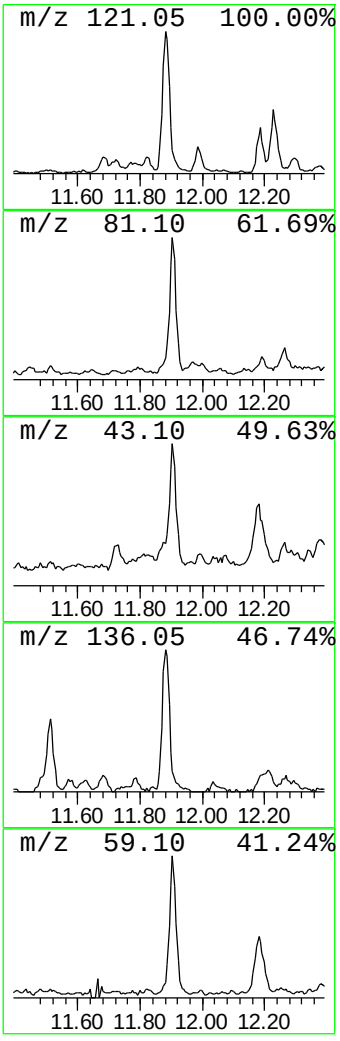
Peak Number 17 Phenol, 3,4,5-trimethyl- Concentration Rank 23

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.89	3.96 ng/ul	263803	Naphthalene-d8	10.14

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, 3,4,5-trimethyl-	136	C9H12O	000527-54-8	92
2	Phenol, 3-(1-methylethyl)-	136	C9H12O	000618-45-1	92
3	Phenol, 2,3,6-trimethyl-	136	C9H12O	002416-94-6	81
4	Phenol, 2,3,5-trimethyl-	136	C9H12O	000697-82-5	81
5	Phenol, 2,3,6-trimethyl-	136	C9H12O	002416-94-6	81





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Acq On : 24 Jun 2020 03:30
Operator : JU/CG
Sample : L3069-12
Misc :
ALS Vial : 29 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : CB7H2

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM061220MA.M
Quant Title : SVOA CALIBRATION

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

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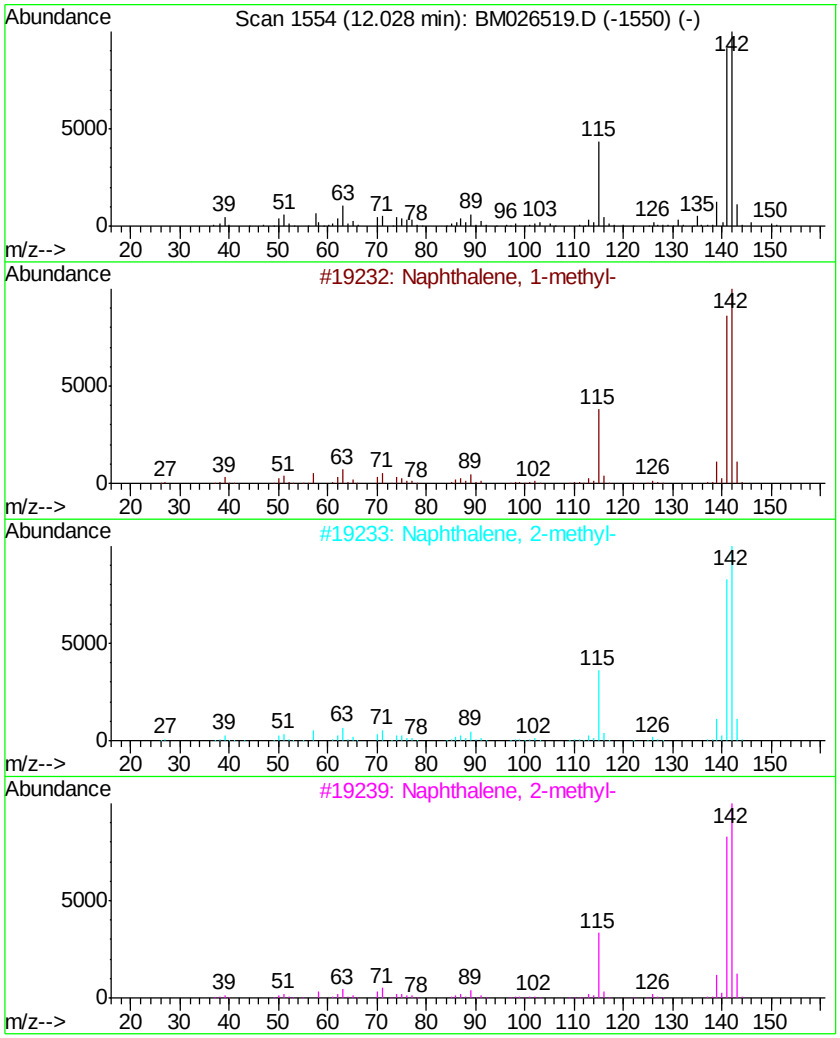
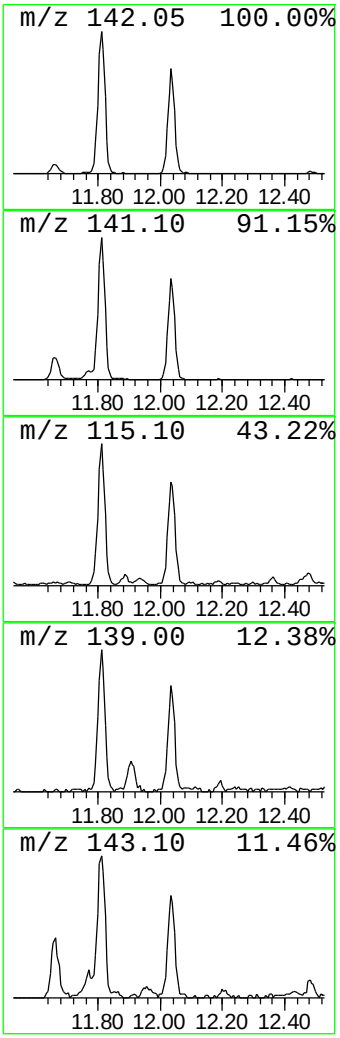
Peak Number 18 Naphthalene, 1-methyl- Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.03	7.40 ng/ul	493130	Naphthalene-d8	10.14

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	95
4	Benzocycloheptatriene	142	C11H10	000264-09-5	93
5	1,4-Methanonaphthalene, 1,4-dihy...	142	C11H10	004453-90-1	93





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Operator : JU/CG
Sample : L3069-12
Misc :
ALS Vial : 29 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : CB7H2

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM061220MA.M
Quant Title : SVOA CALIBRATION

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

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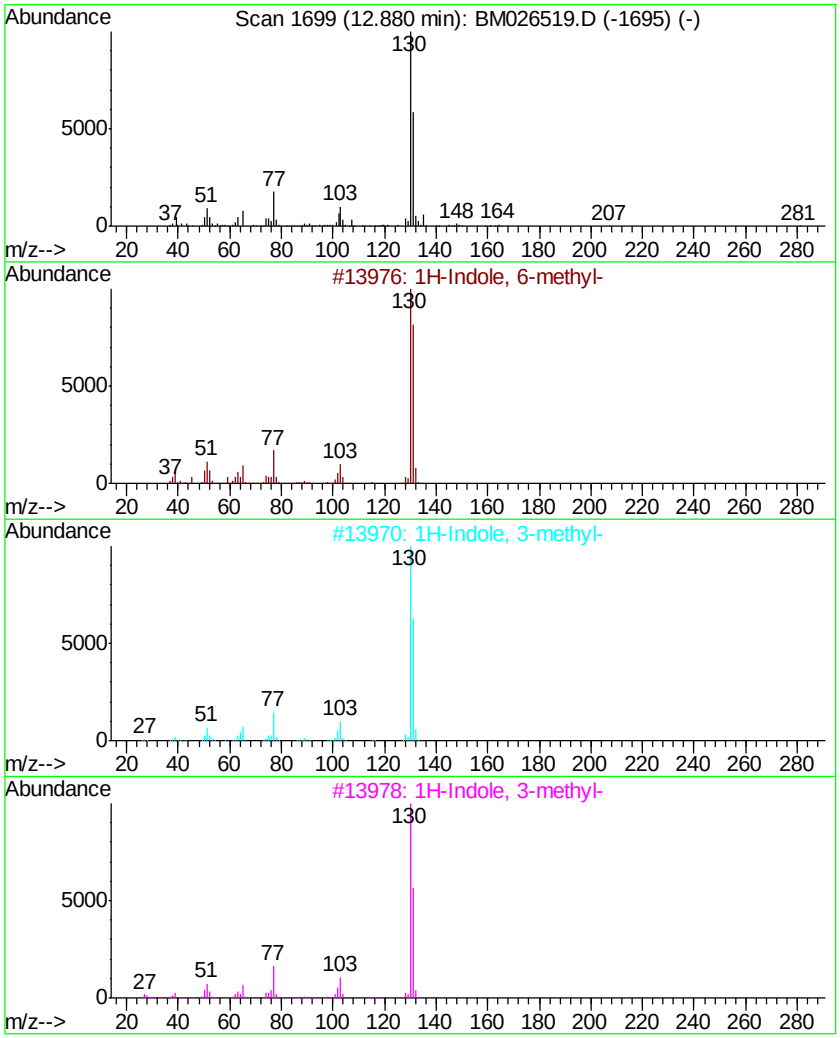
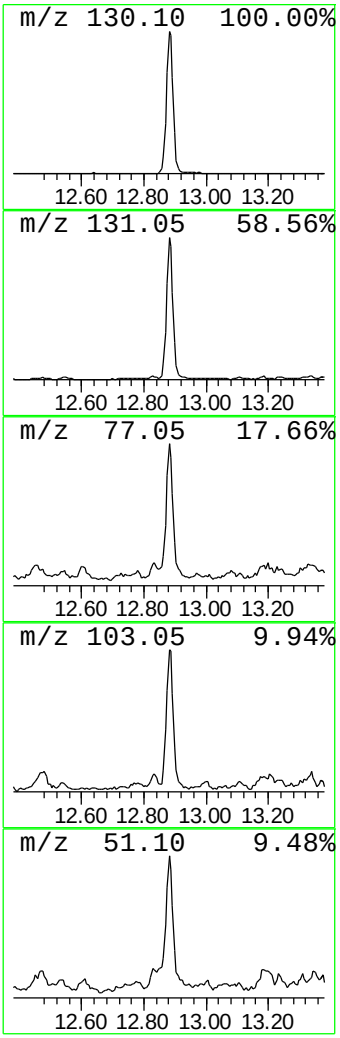
Peak Number 19 1H-Indole, 6-methyl- Concentration Rank 22

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.88	3.97 ng/ul	570959	Acenaphthene-d10	14.04

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Indole, 6-methyl-	131	C9H9N	003420-02-8	97
2	1H-Indole, 3-methyl-	131	C9H9N	000083-34-1	94
3	1H-Indole, 3-methyl-	131	C9H9N	000083-34-1	94
4	1H-Indole, 7-methyl-	131	C9H9N	000933-67-5	92
5	1H-Indole, 3-methyl-	131	C9H9N	000083-34-1	91





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Acq On : 24 Jun 2020 03:30
Operator : JU/CG
Sample : L3069-12
Misc :
ALS Vial : 29 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : CB7H2

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM061220MA.M
Quant Title : SVOA CALIBRATION

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

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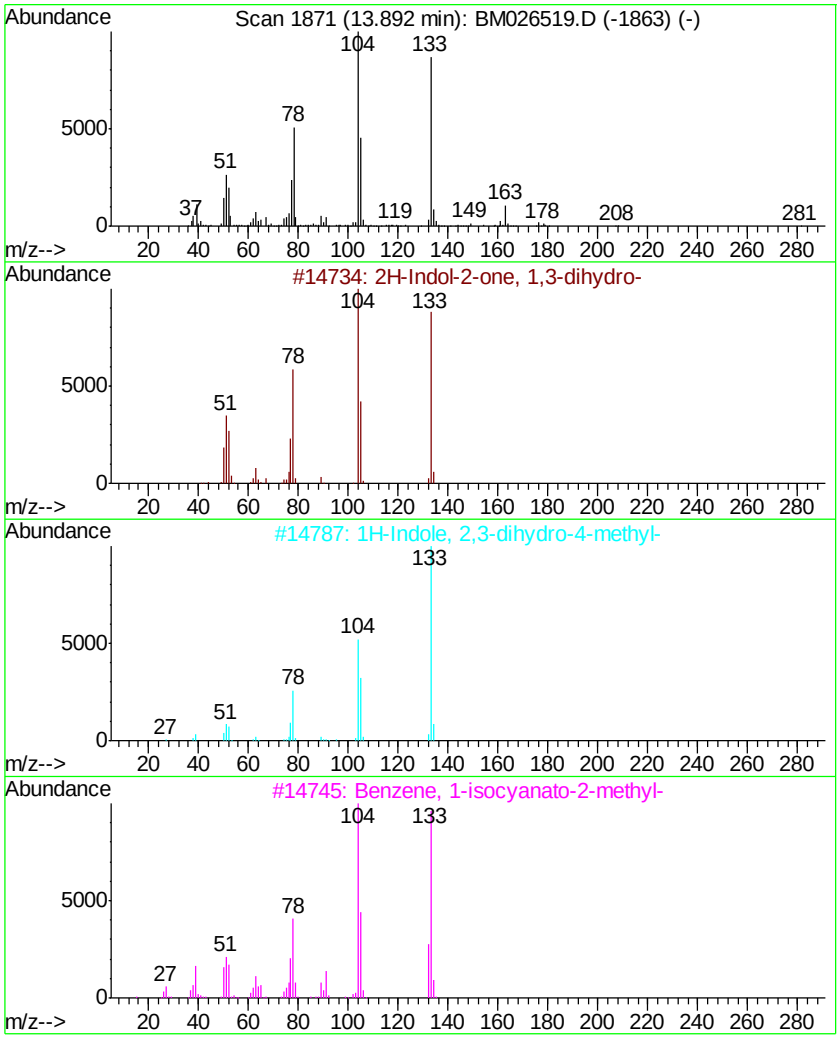
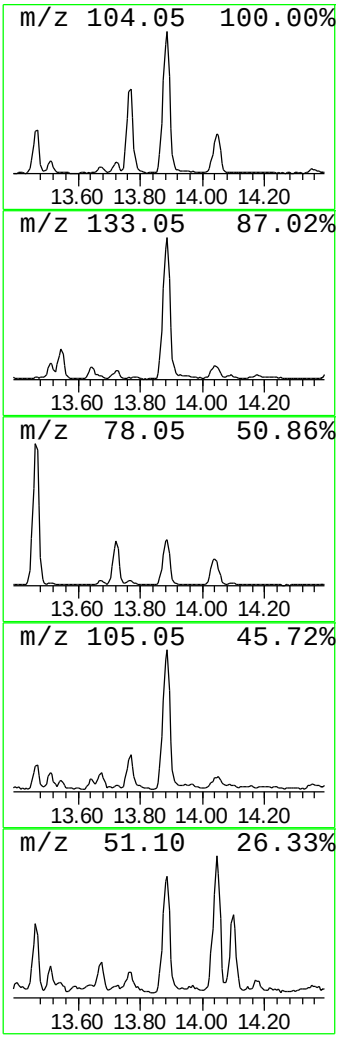
Peak Number 20 2H-Indol-2-one, 1,3-dihydro- Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.89	4.52 ng/ul	649725	Acenaphthene-d10	14.04

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2H-Indol-2-one, 1,3-dihydro-	133	C8H7NO	000059-48-3	97
2	1H-Indole, 2,3-dihydro-4-methyl-	133	C9H11N	062108-16-1	96
3	Benzene, 1-isocyanato-2-methyl-	133	C8H7NO	000614-68-6	91
4	Benzene, 1-isocyanato-2-methyl-	133	C8H7NO	000614-68-6	91
5	1H-Benzotriazole, 4-methyl-	133	C7H7N3	029878-31-7	90





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Operator : JU/CG
Sample : L3069-12
Misc :
ALS Vial : 29 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : CB7H2

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Quant Title : SVOA CALIBRATION

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

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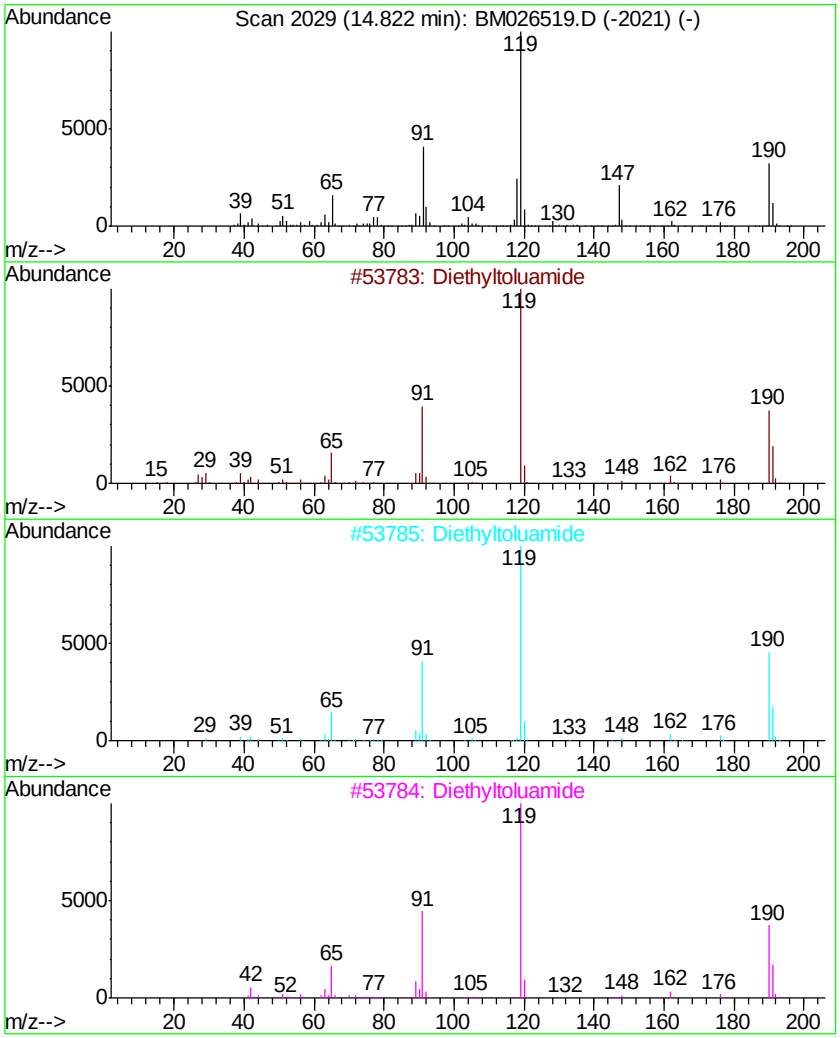
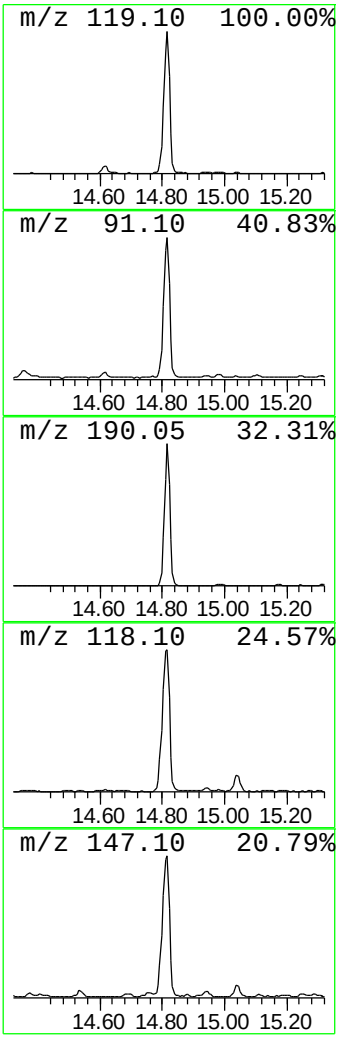
Peak Number 21 Diethyltoluamide Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.82	15.56 ng/ul	2236210	Acenaphthene-d10	14.04

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Diethyltoluamide	191	C12H17NO	000134-62-3	94
2	Diethyltoluamide	191	C12H17NO	000134-62-3	92
3	Diethyltoluamide	191	C12H17NO	000134-62-3	70
4	Benzamide, N,N-diethyl-4-methyl-	191	C12H17NO	002728-05-4	52
5	Benzonitrile, 2-ethoxy-	147	C9H9NO	006609-57-0	52





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Sample : L3069-12
Misc :
ALS Vial : 29 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : CB7H2

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Quant Title : SVOA CALIBRATION

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

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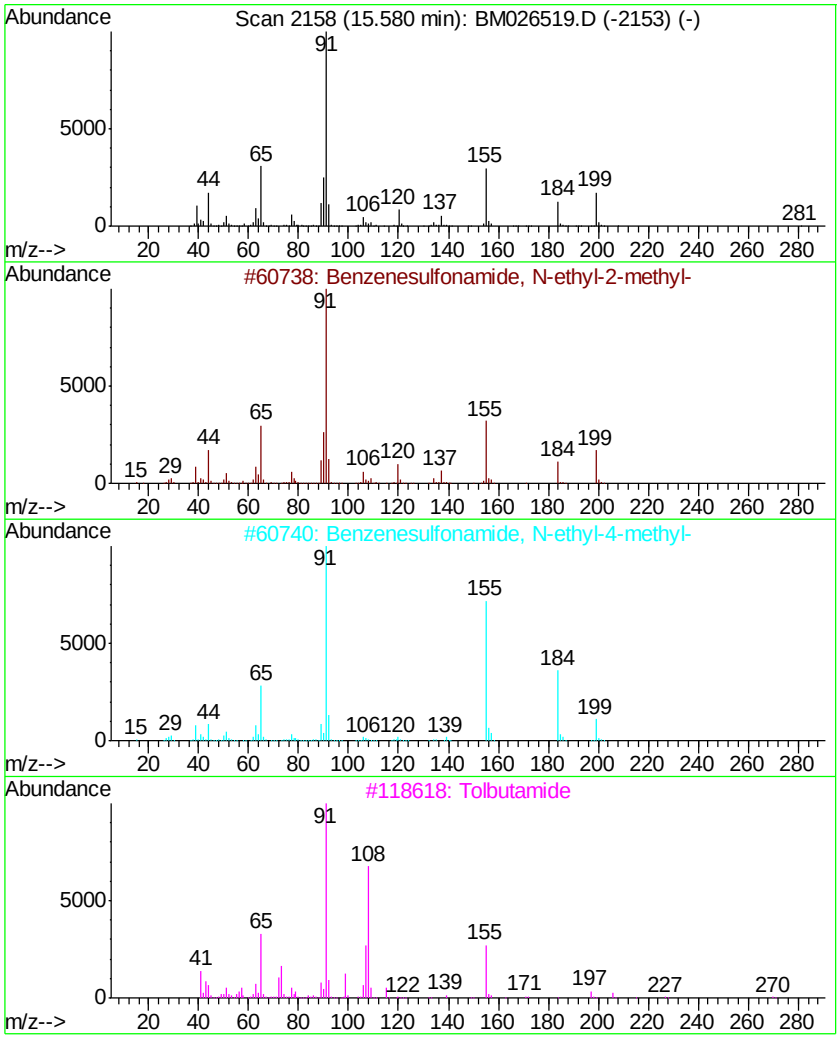
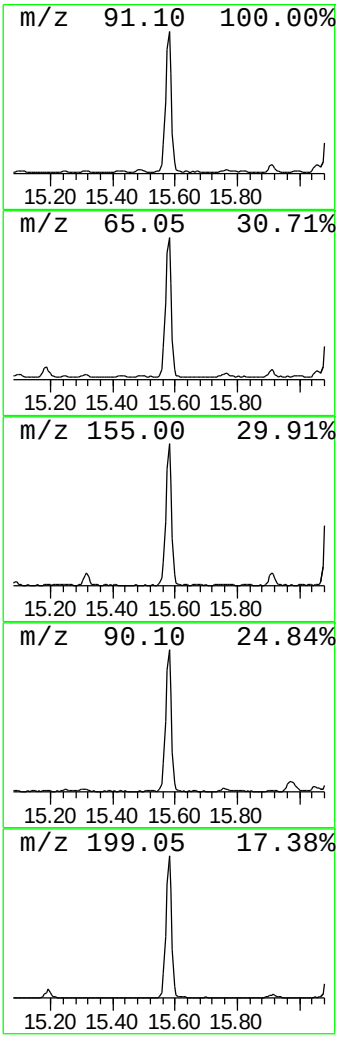
Peak Number 22 Benzenesulfonamide, N-ethyl... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.58	12.45 ng/ul	1376870	Phenanthrene-d10	16.79

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-4-me...	199	C9H13NO2S	000080-39-7	59
3	Tolbutamide	270	C12H18N2O3S	000064-77-7	50
4	p-Toluenesulfonylacetonitrile	195	C9H9NO2S	005697-44-9	50
5	Benzenesulfonamide, N-ethyl-4-me...	199	C9H13NO2S	000080-39-7	47





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Acq On : 24 Jun 2020 03:30
Operator : JU/CG
Sample : L3069-12
Misc :
ALS Vial : 29 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : CB7H2

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Quant Title : SVOA CALIBRATION

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

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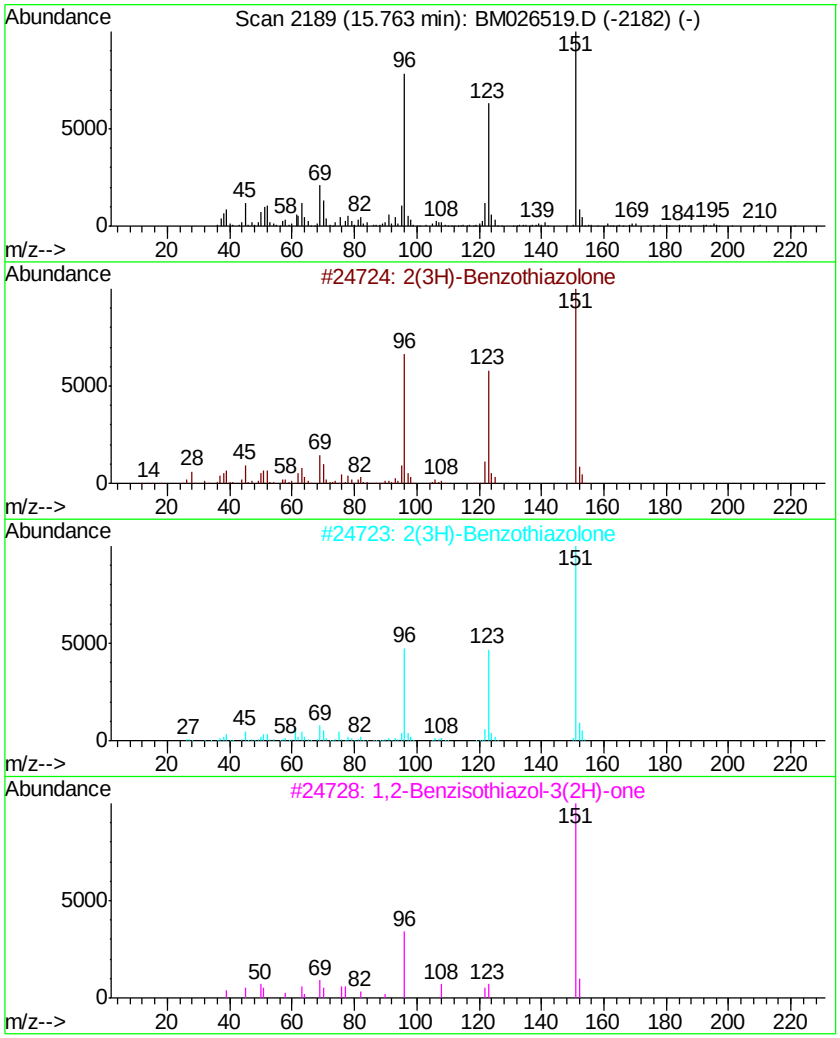
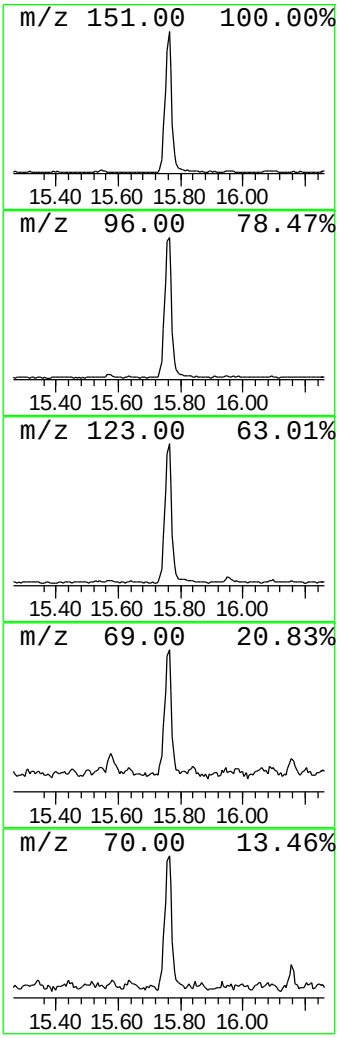
Peak Number 23 2(3H)-Benzothiazolone Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.76	6.04 ng/ul	667968	Phenanthrene-d10	16.79

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	91
3	1,2-Benzisothiazol-3(2H)-one	151	C7H5NOS	002634-33-5	68
4	5-Cyclohexyl-1-pentene	152	C11H20	005729-54-4	38
5	1,2,4-Triazolo[4,3-a]pyridine-3(...	151	C6H5N3S	006952-68-7	27





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Sample : L3069-12
Misc :
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Instrument : BNA_M
ClientSampleId : CB7H2

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Quant Title : SVOA CALIBRATION

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

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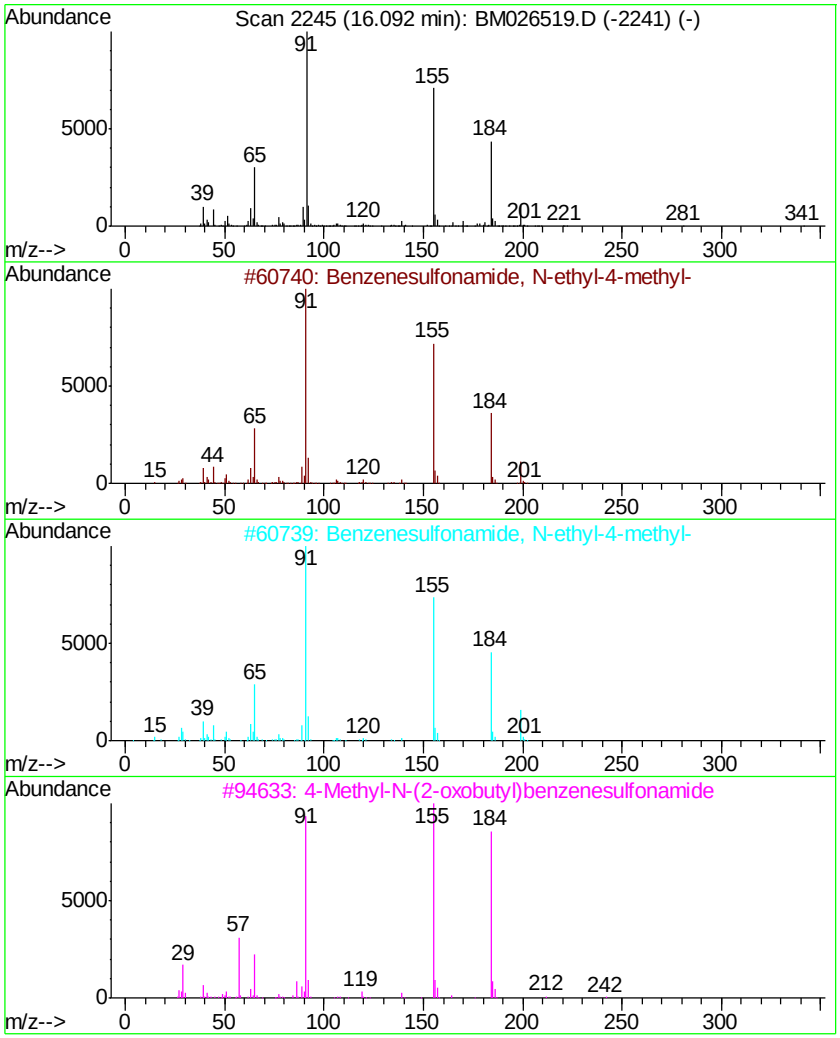
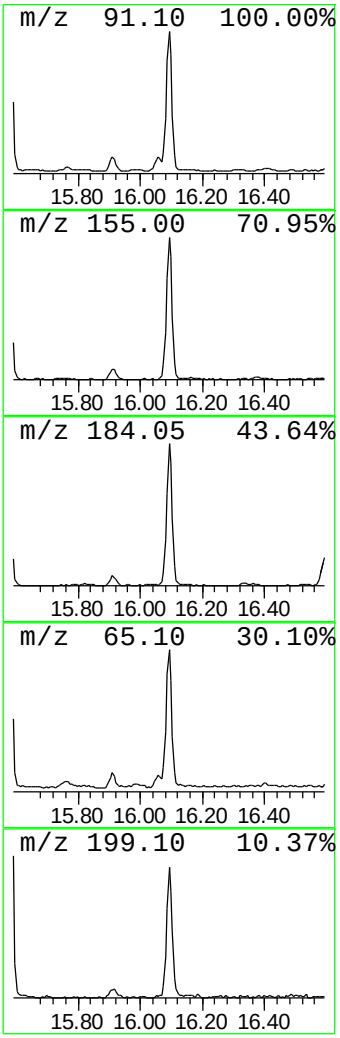
Peak Number 24 Benzenesulfonamide, N-ethyl... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.09	7.74 ng/ul	856127	Phenanthrene-d10	16.79

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzenesulfonamide, N-ethyl-4-me...	199	C9H13NO2S	000080-39-7	92
2	Benzenesulfonamide, N-ethyl-4-me...	199	C9H13NO2S	000080-39-7	87
3	4-Methyl-N-(2-oxobutyl)benzenesu...	241	C11H15NO3S	1000192-14-5	86
4	4-Methyl-N-[2-(3-nitro-[1,2,4]tr...	311	C11H13N5O4S	1000301-03-5	78
5	Furazan-3-carboxylic acid, 4-ami...	325	C12H15N5O4S	1000304-69-4	78





Library Search Compound Report

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Acq On : 24 Jun 2020 03:30
Operator : JU/CG
Sample : L3069-12
Misc :
ALS Vial : 29 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : CB7H2

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM061220MA.M
Quant Title : SVOA CALIBRATION

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

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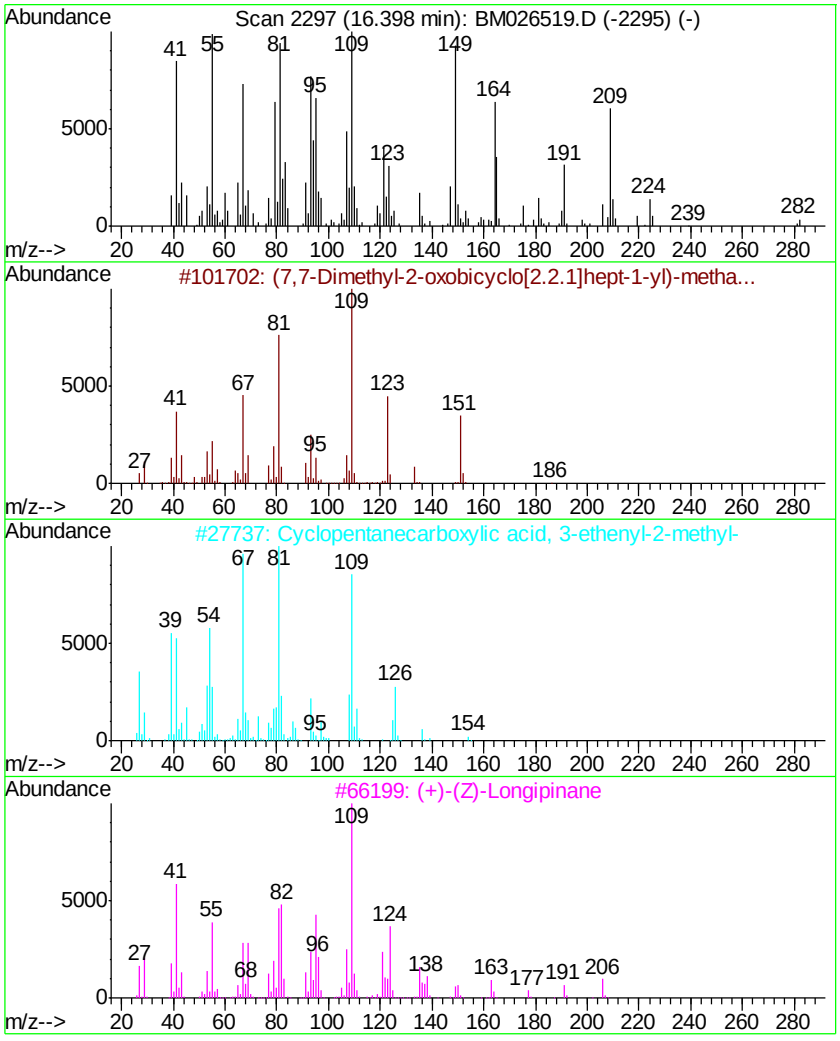
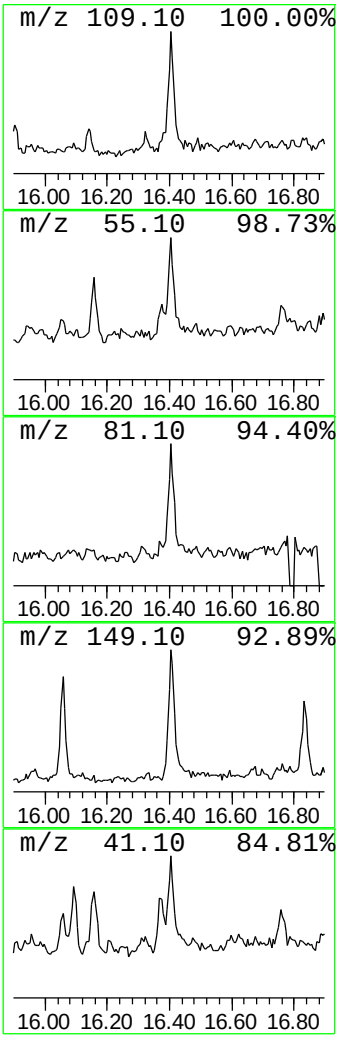
Peak Number 25 unknown-02 Concentration Rank 30

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.40	2.14 ng/ul	237255	Phenanthrene-d10	16.79

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	(7,7-Dimethyl-2-oxobicyclo[2.2.1...	250	C10H15ClO3S	1000194-76-1	27
2	Cyclopentanecarboxylic acid, 3-e...	154	C9H14O2	108451-44-1	22
3	(+)-(Z)-Longipinane	206	C15H26	1000156-12-3	20
4	Cyclohexene, 3,3,5-trimethyl-	124	C9H16	000503-45-7	20
5	1H-Indene, octahydro-2,3a,4-trim...	208	C15H28	031230-13-4	20





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Misc :
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Instrument : BNA_M
ClientSampleId : CB7H2

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Quant Title : SVOA CALIBRATION

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

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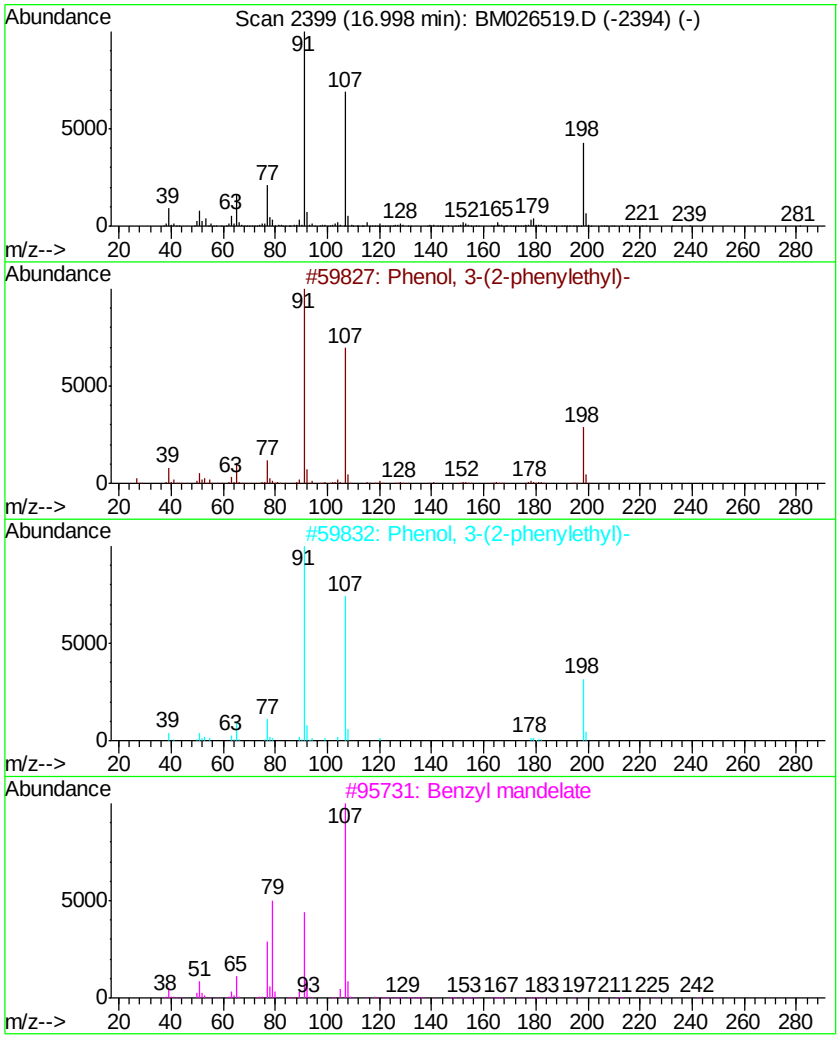
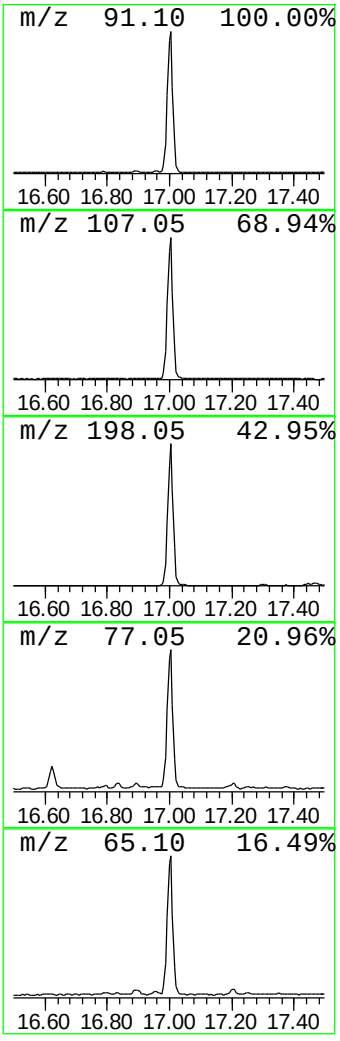
Peak Number 26 Phenol, 3-(2-phenylethyl)- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.00	22.90 ng/ul	2533540	Phenanthrene-d10	16.79

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, 3-(2-phenylethyl)-	198	C14H14O	033675-75-1	93
2	Phenol, 3-(2-phenylethyl)-	198	C14H14O	033675-75-1	93
3	Benzyl mandelate	242	C15H14O3	000890-98-2	43
4	Benzenemethanol, .alpha.-pentyl-	178	C12H18O	004471-05-0	38
5	2-Phenyl-4-(cyanomethyl)-5-methy...	198	C11H10N4	013322-20-8	38





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Instrument : BNA_M
ClientSampleId : CB7H2

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Quant Title : SVOA CALIBRATION

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

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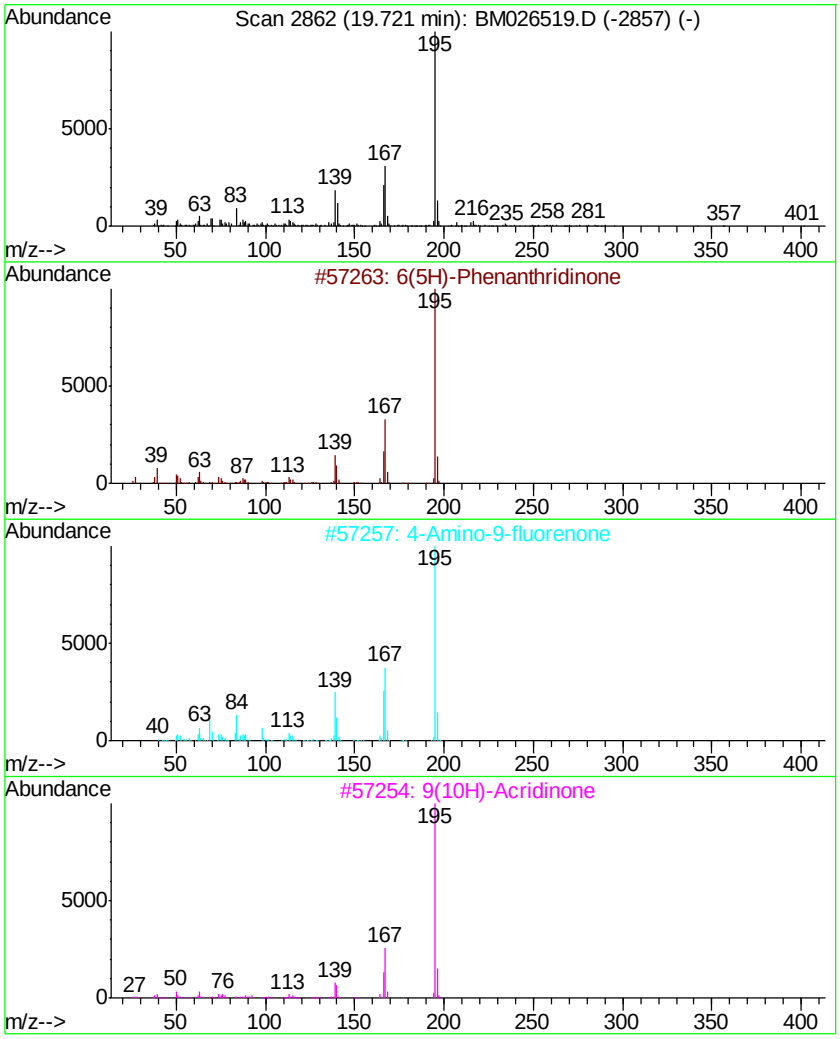
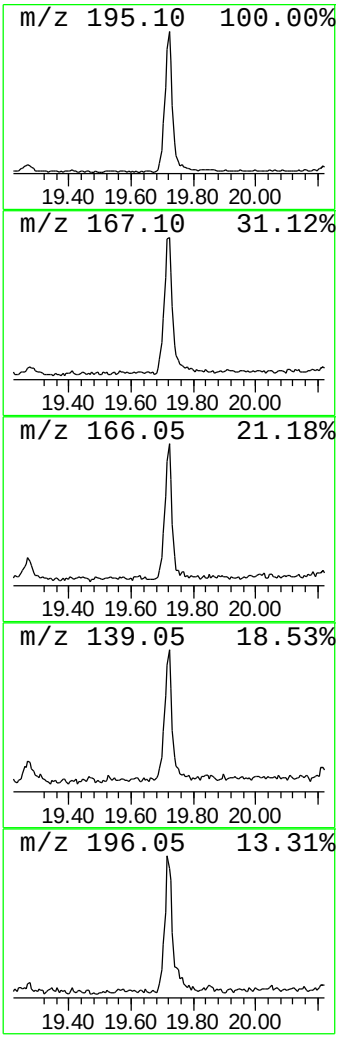
Peak Number 27 6(5H)-Phenanthridinone Concentration Rank 27

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.72	2.70 ng/ul	323339	Chrysene-d12	21.00

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	6(5H)-Phenanthridinone	195	C13H9NO	001015-89-0	95
2	4-Amino-9-fluorenone	195	C13H9NO	004269-15-2	94
3	9(10H)-Acridinone	195	C13H9NO	000578-95-0	94
4	4-Amino-9-fluorenone	195	C13H9NO	004269-15-2	94
5	9(10H)-Acridinone	195	C13H9NO	000578-95-0	94





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Sample : L3069-12
Misc :
ALS Vial : 29 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : CB7H2

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM061220MA.M
Quant Title : SVOA CALIBRATION

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

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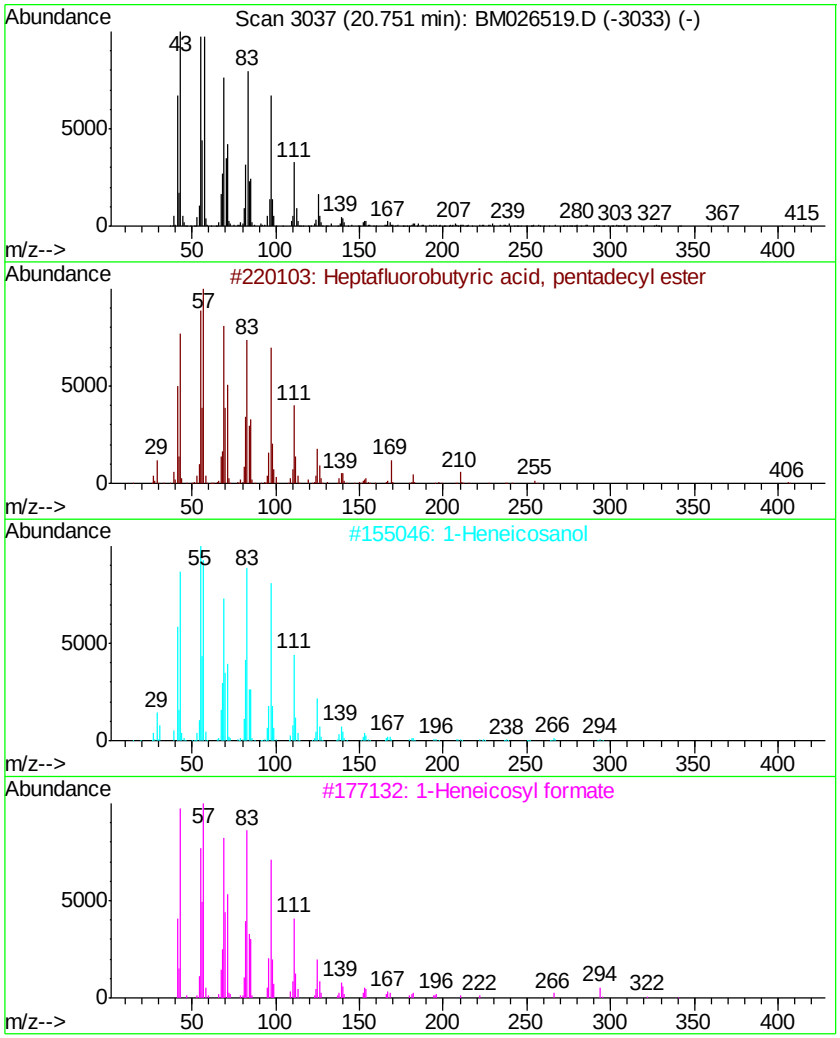
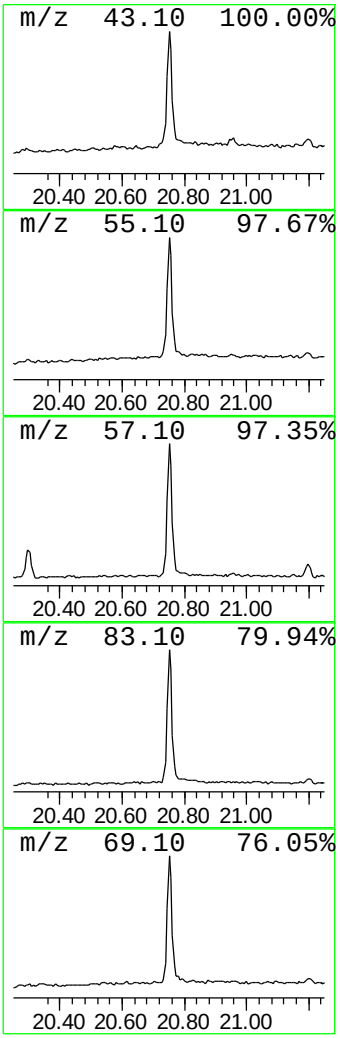
Peak Number 28 Heptafluorobutyric acid, pe... Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.75	7.43 ng/ul	889978	Chrysene-d12	21.00

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptafluorobutyric acid, pentade...	424	C19H31F7O2	959261-23-5	94
2	1-Heneicosanol	312	C21H44O	015594-90-8	94
3	1-Heneicosyl formate	340	C22H44O2	077899-03-7	94
4	Pentafluoropropionic acid, penta...	374	C18H31F5O2	959092-08-1	91
5	n-Tetracosanol-1	354	C24H50O	000506-51-4	91





Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM062320\
Data File : BM026519.D
Acq On : 24 Jun 2020 03:30
Operator : JU/CG
Sample : L3069-12
Misc :
ALS Vial : 29 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : CB7H2

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM061220MA.M
Quant Title : SVOA CALIBRATION

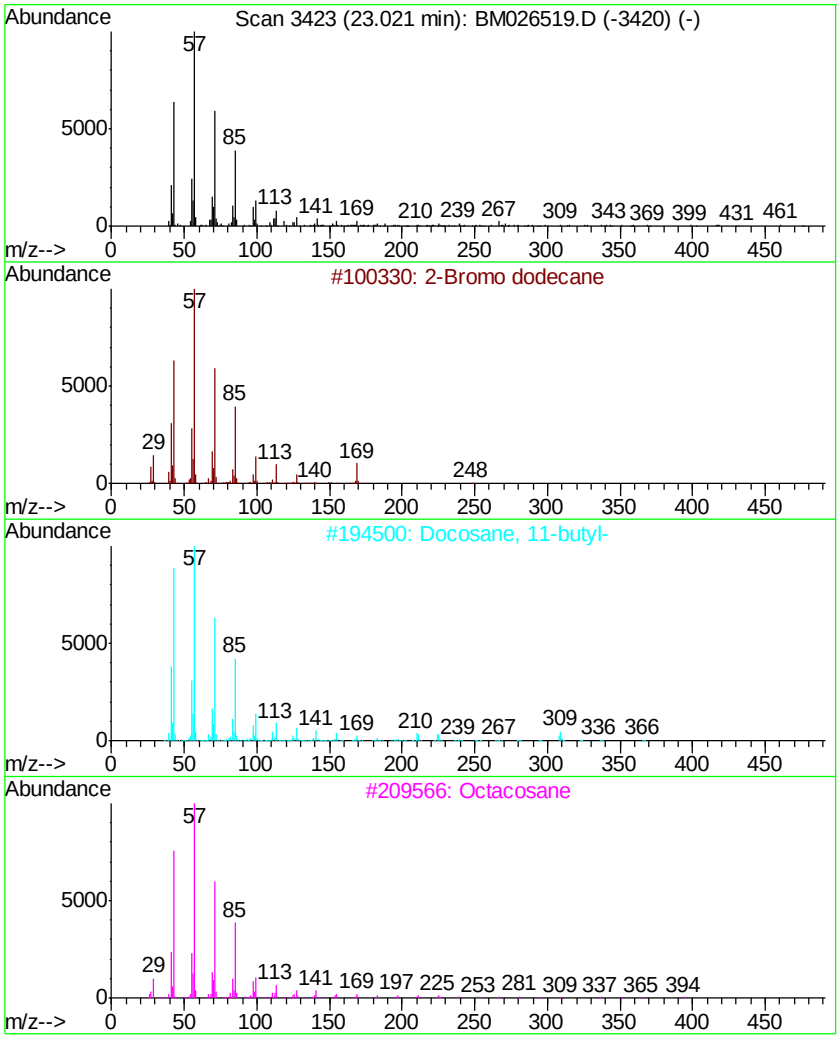
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

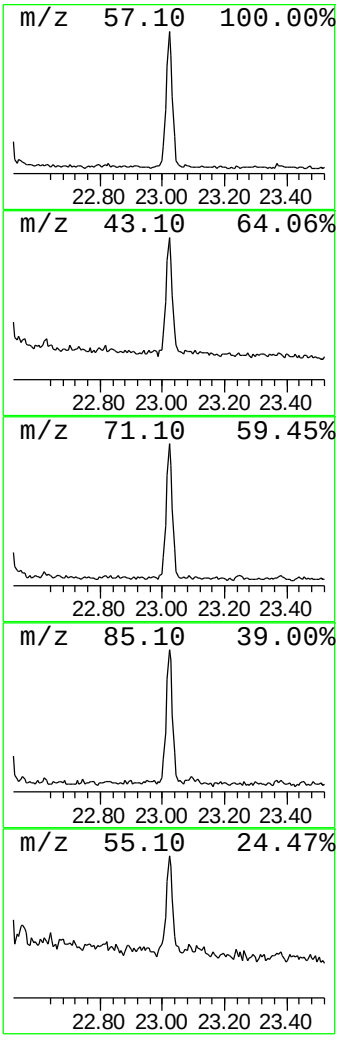
****

Peak Number 29 2-Bromo dodecane Concentration Rank 28

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.02	2.62 ng/ul	242936	Perylene-d12	23.09

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Bromo dodecane	248	C12H25Br	013187-99-0	91
2	Docosane, 11-butyl-	366	C26H54	013475-76-8	90
3	Octacosane	394	C28H58	000630-02-4	90
4	Hexacosane	366	C26H54	000630-01-3	87
5	Tetracosane	338	C24H50	000646-31-1	87





Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM062320\
Data File : BM026519.D
Acq On : 24 Jun 2020 03:30
Operator : JU/CG
Sample : L3069-12
Misc :
ALS Vial : 29 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : CB7H2

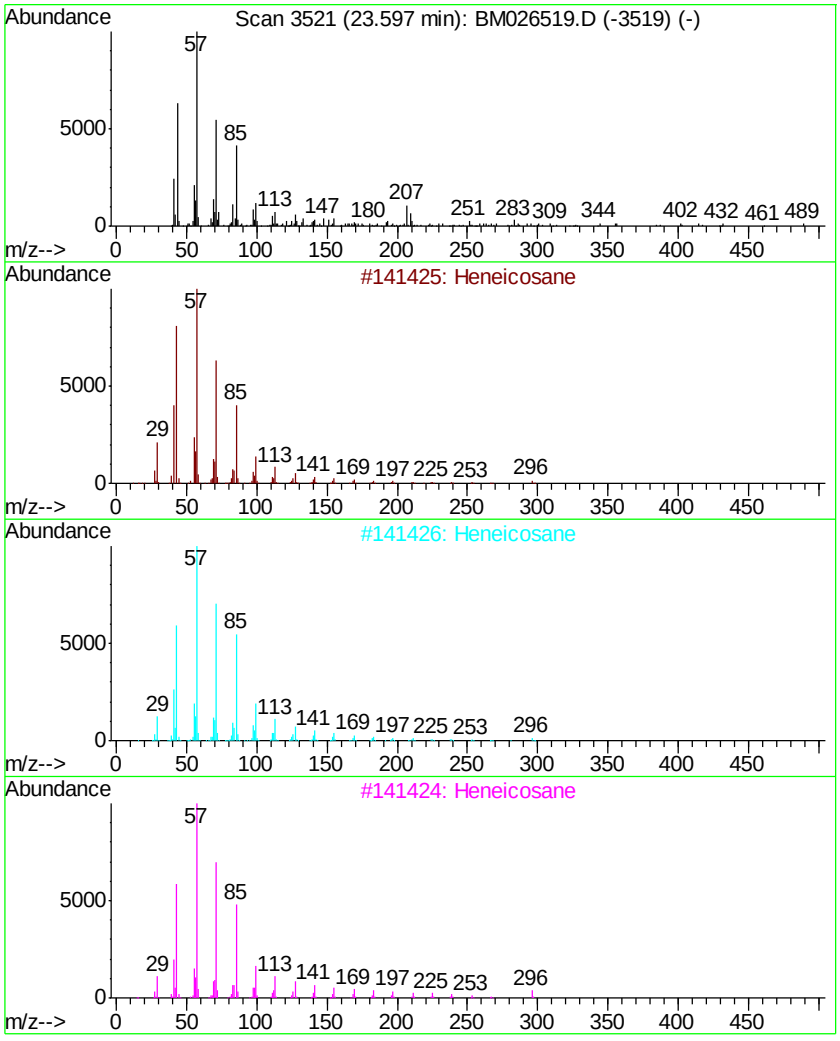
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM061220MA.M
Quant Title : SVOA CALIBRATION

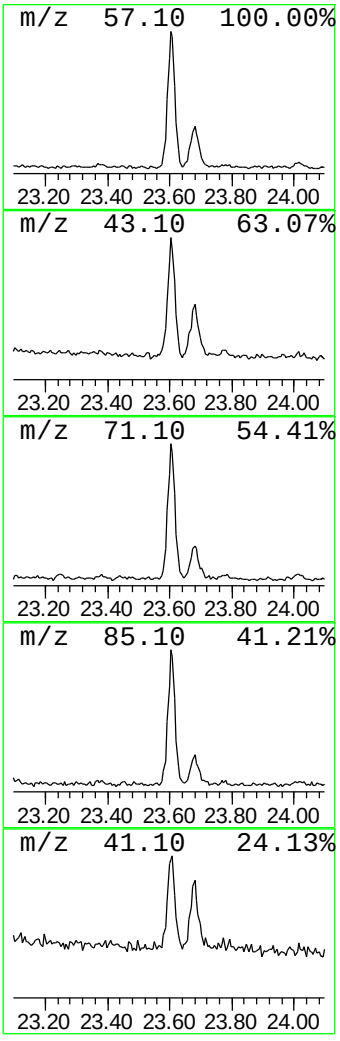
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 30 (DEL) Alkane: Straight-Chai... Concentration Rank 26

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.60	2.73 ng/ul	253308	Perylene-d12	23.09

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heneicosane	296	C21H44	000629-94-7	97
2	Heneicosane	296	C21H44	000629-94-7	94
3	Heneicosane	296	C21H44	000629-94-7	87
4	Tetradecane, 2,6,10-trimethyl-	240	C17H36	014905-56-7	87
5	Octadecane	254	C18H38	000593-45-3	83





Tentatively Identified Compound (LSC) summary

Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM062320\
Data File : BM026519.D
Acq On : 24 Jun 2020 03:30
Operator : JU/CG
Sample : L3069-12
Misc :
ALS Vial : 29 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : CB7H2

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM061220MA.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
Toluene	3.66	4.5	ng/ul	196162	1	7.38	875923	20.0
1,3-Oxathiolane	4.09	3.1	ng/ul	134463	1	7.38	875923	20.0
p-Cymene	7.52	75.1	ng/ul	3290950	1	7.38	875923	20.0
Indane	7.74	4.9	ng/ul	214424	1	7.38	875923	20.0
Benzenamine, 3-me...	8.36	8.6	ng/ul	376644	1	7.38	875923	20.0
Bicyclo[2.2.1]hep...	8.61	30.2	ng/ul	1323720	1	7.38	875923	20.0
Cyclohexanol, 1-m...	9.39	4.8	ng/ul	323477	2	10.14	1333060	20.0
Cyclohexanemethan...	9.53	60.0	ng/ul	4001950	2	10.14	1333060	20.0
Camphor	9.56	44.5	ng/ul	2967940	2	10.14	1333060	20.0
unknown-01	9.69	10.0	ng/ul	667926	2	10.14	1333060	20.0
1,4-Pentadiene, 2...	9.92	2.5	ng/ul	167410	2	10.14	1333060	20.0
Terpinen-4-ol	10.02	22.4	ng/ul	1495700	2	10.14	1333060	20.0
Bicyclo[3.1.1]hep...	10.50	3.2	ng/ul	213628	2	10.14	1333060	20.0
Phenol, 2-propyl-	10.99	9.6	ng/ul	636640	2	10.14	1333060	20.0
Phenol, p-tert-bu...	11.51	4.5	ng/ul	302528	2	10.14	1333060	20.0
Indole	11.66	4.3	ng/ul	289030	2	10.14	1333060	20.0
Phenol, 3,4,5-tri...	11.89	4.0	ng/ul	263803	2	10.14	1333060	20.0
Naphthalene, 1-me...	12.03	7.4	ng/ul	493130	2	10.14	1333060	20.0
1H-Indole, 6-methyl-	12.88	4.0	ng/ul	570959	3	14.04	2874140	20.0
2H-Indol-2-one, 1...	13.89	4.5	ng/ul	649725	3	14.04	2874140	20.0
Diethyltoluamide	14.82	15.6	ng/ul	2236210	3	14.04	2874140	20.0
Benzenesulfonamid...	15.58	12.4	ng/ul	1376870	4	16.79	2212240	20.0
2(3H)-Benzothiazo...	15.76	6.0	ng/ul	667968	4	16.79	2212240	20.0
Benzenesulfonamid...	16.09	7.7	ng/ul	856127	4	16.79	2212240	20.0
unknown-02	16.40	2.1	ng/ul	237255	4	16.79	2212240	20.0
Phenol, 3-(2-phen...	17.00	22.9	ng/ul	2533540	4	16.79	2212240	20.0
6(5H)-Phenanthrid...	19.72	2.7	ng/ul	323339	5	21.00	2395450	20.0
Heptafluorobutyri...	20.75	7.4	ng/ul	889978	5	21.00	2395450	20.0
2-Bromo dodecane	23.02	2.6	ng/ul	242936	6	23.09	1852860	20.0
(DEL) Alkane: Str...	23.60	2.7	ng/ul	253308	6	23.09	1852860	20.0