LSC Area Percent Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM052920\
Data File : BM026187.D
Acq On : 30 May 2020 00:46
Operator : MA/SJ
Sample : L2820-08
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : CB638

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	3.052	26	28	30	rBV	46141	48856	0.64%	0.068%
2	3.081	30	33	48	rVB	507234	700039	9.12%	0.980%
3	3.840	158	162	165	rBV	101041	122966	1.60%	0.172%
4	3.875	165	168	174	rVB	28069	36603	0.48%	0.051%
5	4.181	215	220	233	rVV	176883	280910	3.66%	0.393%
6	4.840	328	332	338	rBV	28207	44538	0.58%	0.062%
7	5.163	383	387	396	rVB3	24771	51343	0.67%	0.072%
8	5.793	487	494	499	rBV	26045	46321	0.60%	0.065%
9	6.181	555	560	569	rVV	60431	100999	1.32%	0.141%
10	6.481	598	611	616	rBV2	110059	234027	3.05%	0.328%
11	6.669	637	643	657	rVV	149578	277086	3.61%	0.388%
12	6.828	664	670	677	rBV	387665	618611	8.06%	0.866%
13	7.016	695	702	710	rVV	481228	776014	10.11%	1.087%
14	7.475	771	780	788	rBV	572142	941387	12.26%	1.318%
15	7.551	788	793	797	rVV2	48408	81624	1.06%	0.114%
16	7.616	797	804	813	rVV	2330638	3398970	44.28%	4.759%
17	7.781	826	832	838	rVV2	55495	91302	1.19%	0.128%
18	7.840	838	842	849	rVB2	31864	52121	0.68%	0.073%
19	8.193	896	902	909	rVV	416757	653747	8.52%	0.915%
20	8.251	909	912	920	rVB6	22351	43555	0.57%	0.061%
21	8.363	925	931	936	rBV4	27436	62047	0.81%	0.087%
22	8.457	940	947	954	rBV	5007000	7676538	100.00%	10.748%
23	8.575	962	967	970	rBV2	50730	82255	1.07%	0.115%
24	8.616	970	974	983	rVV	496691	820695	10.69%	1.149%
25	8.704	983	989	996	rVB	1332977	1974190	25.72%	2.764%
26	8.822	1003	1009	1011	rBV4	24478	41509	0.54%	0.058%
27	8.857	1011	1015	1028	rVB8	41582	88068	1.15%	0.123%
28	8.975	1028	1035	1041	rBV3	48865	97073	1.26%	0.136%
29	9.104	1050	1057	1063	rBV	115783	216906	2.83%	0.304%
30	9.234	1075	1079	1082	rBV3	32862	48236	0.63%	0.068%
31	9.334	1090	1096	1101	rBV	424180	694377	9.05%	0.972%
32	9.381	1102	1104	1111	rVB2	54394	68301	0.89%	0.096%
33	9.463	1111	1118	1121	rBV6	45808	92248	1.20%	0.129%
34	9.510	1121	1126	1132	rVV4	89377	187654	2.44%	0.263%

LSC Area Percent Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM052920\
Data File : BM026187.D
Acq On : 30 May 2020 00:46
Operator : MA/SJ
Sample : L2820-08
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : CB638

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

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

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

35	9.622	1135	1145	1149	rVV	3088428	5755920	74.98%	8.059%
36	9.663	1149	1152	1161	rVV	2254734	3259614	42.46%	4.564%
37	9.792	1167	1174	1180	rVB	219126	385074	5.02%	0.539%
38	9.869	1180	1187	1192	rBV	812840	1364560	17.78%	1.911%
39	9.916	1192	1195	1205	rVB2	145841	283526	3.69%	0.397%
40	10.034	1208	1215	1222	rBV5	43212	105538	1.37%	0.148%
41	10.116	1222	1229	1235	rBV	401467	629469	8.20%	0.881%
42	10.234	1243	1249	1255	rVB	753288	1190080	15.50%	1.666%
43	10.304	1256	1261	1266	rVB3	122586	199206	2.59%	0.279%
44	10.381	1266	1274	1286	rVB	758661	1401171	18.25%	1.962%
45	10.504	1289	1295	1297	rBV6	29724	58390	0.76%	0.082%
46	10.598	1307	1311	1315	rVB4	30646	45551	0.59%	0.064%
47	10.698	1323	1328	1331	rBV6	28182	45969	0.60%	0.064%
48	10.828	1345	1350	1355	rVV7	33453	68856	0.90%	0.096%
49	10.886	1356	1360	1363	rVV5	37676	60505	0.79%	0.085%
50	10.939	1363	1369	1374	rVB5	44074	90399	1.18%	0.127%
51	11.081	1381	1393	1405	rBV	606359	1109228	14.45%	1.553%
52	11.233	1413	1419	1426	rVB9	40594	97958	1.28%	0.137%
53	11.304	1427	1431	1436	rVB5	28515	48706	0.63%	0.068%
54	11.445	1450	1455	1464	rVB6	48176	106847	1.39%	0.150%
55	11.604	1474	1482	1487	rBV	130019	227256	2.96%	0.318%
56	11.698	1494	1498	1501	rBV5	19859	34478	0.45%	0.048%
57	11.751	1501	1507	1514	rBV	1769789	2756762	35.91%	3.860%
58	11.857	1515	1525	1531	rVV	309829	543768	7.08%	0.761%
59	11.992	1540	1548	1555	rVB4	88416	185857	2.42%	0.260%
60	12.092	1561	1565	1568	rVB5	21455	33599	0.44%	0.047%
61	12.269	1590	1595	1599	rBV	42943	71182	0.93%	0.100%
62	12.833	1686	1691	1698	rVV6	66632	132368	1.72%	0.185%
63	12.916	1700	1705	1710	rVV6	43616	101266	1.32%	0.142%
64	12.969	1710	1714	1722	rVB2	67532	113215	1.47%	0.159%
65	13.045	1722	1727	1732	rBV5	32844	70498	0.92%	0.099%
66	13.163	1743	1747	1755	rVB3	55590	101360	1.32%	0.142%
67	13.233	1755	1759	1762	rBV4	25281	42465	0.55%	0.059%
68	13.545	1806	1812	1816	rVV	1194408	1623853	21.15%	2.274%
69	13.592	1816	1820	1824	rVB	207382	266367	3.47%	0.373%
70	13.680	1830	1835	1839	rBV6	41994	78266	1.02%	0.110%
71	13.804	1850	1856	1863	rBV	1196429	1823891	23.76%	2.554%

LSC Area Percent Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM052920\
Data File : BM026187.D
Acq On : 30 May 2020 00:46
Operator : MA/SJ
Sample : L2820-08
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : CB638

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

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

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

72	13.951	1877	1881	1886	rBV	93152	141762	1.85%	0.198%
73	14.045	1893	1897	1904	rVB	136886	193817	2.52%	0.271%
74	14.122	1904	1910	1917	rBV	1149405	1672804	21.79%	2.342%
75	14.357	1945	1950	1955	rBV	103767	168363	2.19%	0.236%
76	14.827	2025	2030	2035	rBV	61650	111851	1.46%	0.157%
77	14.892	2037	2041	2049	rVV2	65493	143141	1.86%	0.200%
78	15.121	2074	2080	2086	rVB	1618920	2197368	28.62%	3.077%
79	15.251	2097	2102	2107	rBV	383705	523975	6.83%	0.734%
80	15.380	2119	2124	2128	rBV3	103674	171830	2.24%	0.241%
81	15.421	2129	2131	2136	rVB5	48315	62966	0.82%	0.088%
82	15.651	2164	2170	2178	rVB	1081708	1478908	19.27%	2.071%
83	15.821	2193	2199	2204	rBV2	344805	549868	7.16%	0.770%
84	15.980	2223	2226	2231	rVB	62947	70762	0.92%	0.099%
85	16.163	2251	2257	2264	rBV	665718	941565	12.27%	1.318%
86	16.221	2265	2267	2271	rVB2	32628	35175	0.46%	0.049%
87	16.439	2301	2304	2306	rBV	74103	85533	1.11%	0.120%
88	16.480	2306	2311	2321	rVB2	700857	995673	12.97%	1.394%
89	16.868	2372	2377	2388	rVB	1492329	2119514	27.61%	2.968%
90	16.968	2388	2394	2404	rBV	1835645	2523955	32.88%	3.534%
91	17.068	2407	2411	2417	rVB	330496	451034	5.88%	0.632%
92	17.804	2532	2536	2542	rBV6	54437	103293	1.35%	0.145%
93	18.074	2578	2582	2586	rBV2	79779	102839	1.34%	0.144%
94	18.562	2661	2665	2676	rVB2	170206	266574	3.47%	0.373%
95	19.268	2780	2785	2792	rBV	2136910	3022590	39.37%	4.232%
96	19.339	2793	2797	2803	rVB	441954	563912	7.35%	0.790%
97	20.821	3046	3049	3055	rVB	438356	483984	6.30%	0.678%
98	21.068	3087	3091	3101	rBV	1887357	2428889	31.64%	3.401%
99	23.056	3423	3429	3436	rVB	1231410	2062296	26.86%	2.888%
100	23.186	3446	3451	3461	rVB	1420656	2482157	32.33%	3.475%

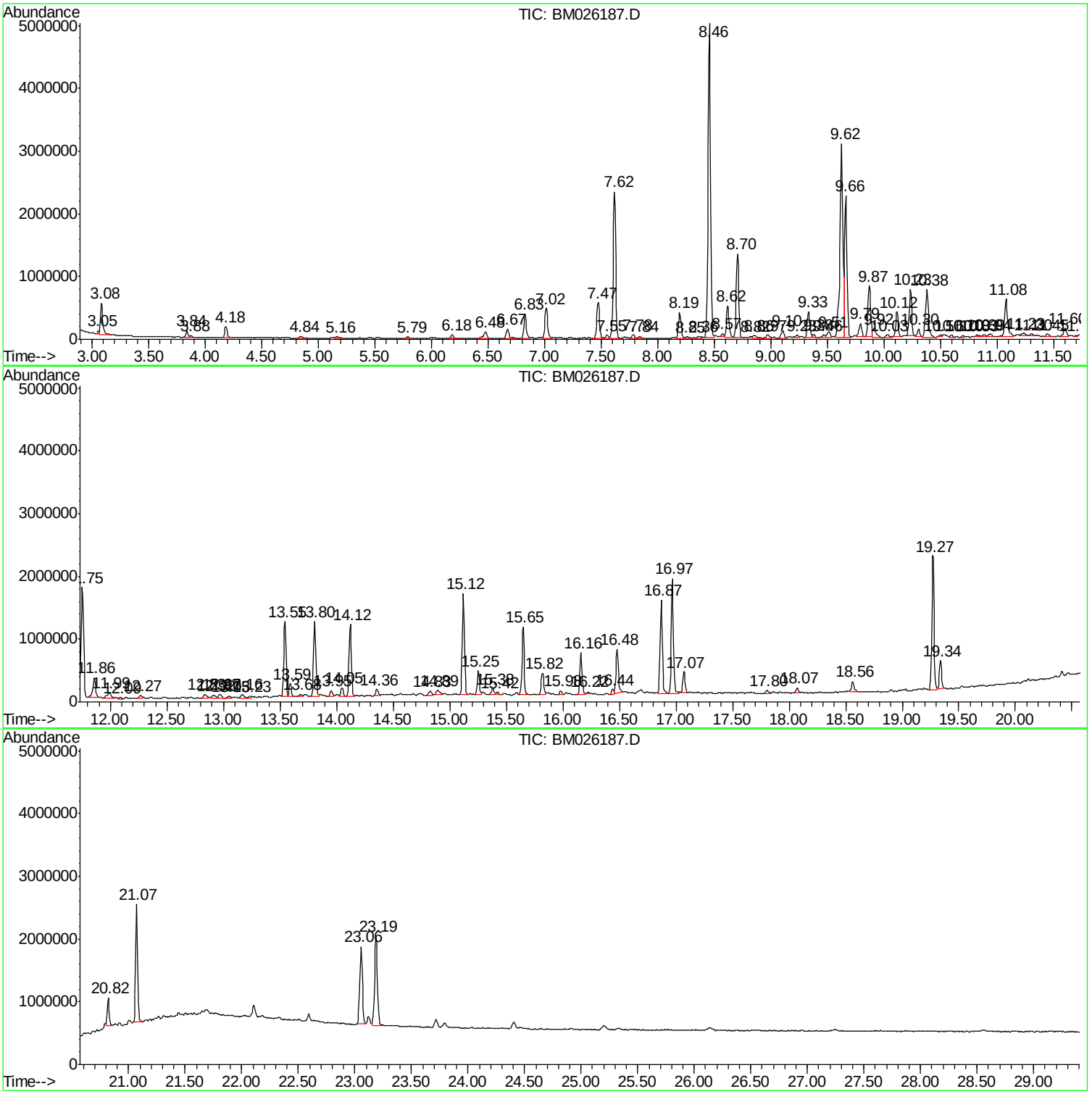
Sum of corrected areas: 71420532

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM052920\
Data File : BM026187.D
Acq On : 30 May 2020 00:46
Operator : MA/SJ
Sample : L2820-08
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
CB638

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 : BM026187.D
Acq On : 30 May 2020 00:46
Operator : MA/SJ
Sample : L2820-08
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : CB638

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

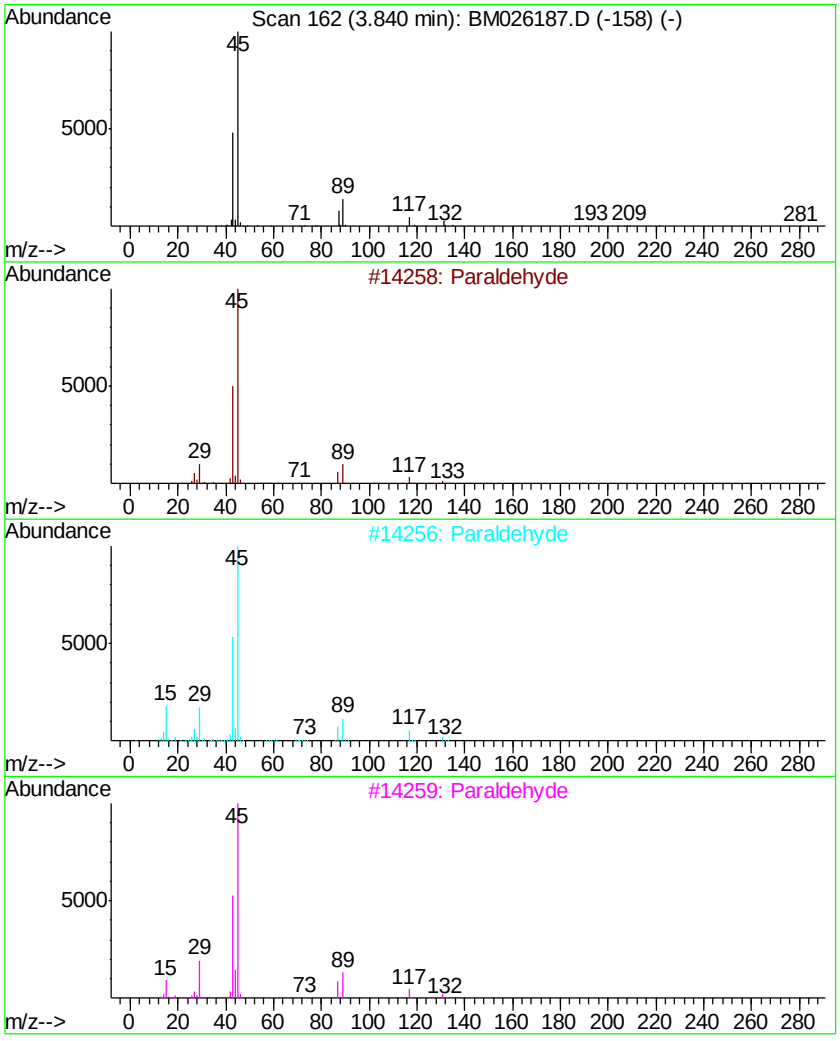
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

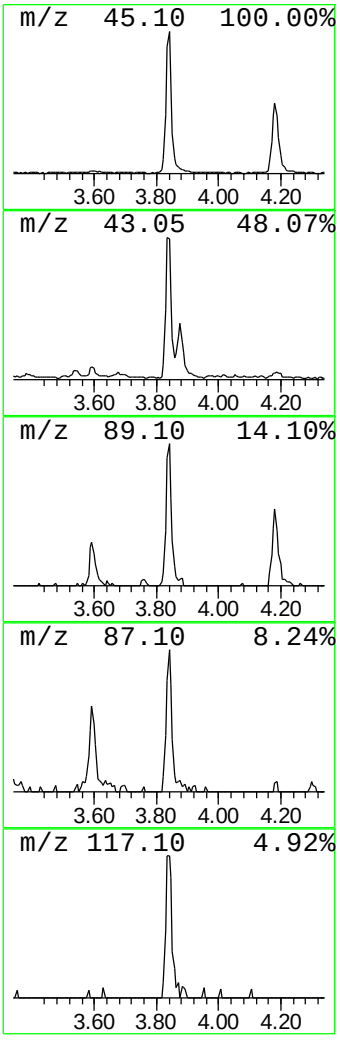
****

Peak Number 1 Paraldehyde Concentration Rank 25

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.84	2.61 ng/ul	122966	1,4-Dichlorobenzene-d4	7.47

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Paraldehyde	132	C6H12O3	000123-63-7	78
2	Paraldehyde	132	C6H12O3	000123-63-7	64
3	Paraldehyde	132	C6H12O3	000123-63-7	64
4	Propane, 2,2'-[ethylidenebis(oxy...	146	C8H18O2	004285-59-0	50
5	Ethanol, 2-[2-(2-butoxyethoxy)et...	206	C10H22O4	000143-22-6	39





Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM052920\
Data File : BM026187.D
Acq On : 30 May 2020 00:46
Operator : MA/SJ
Sample : L2820-08
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : CB638

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

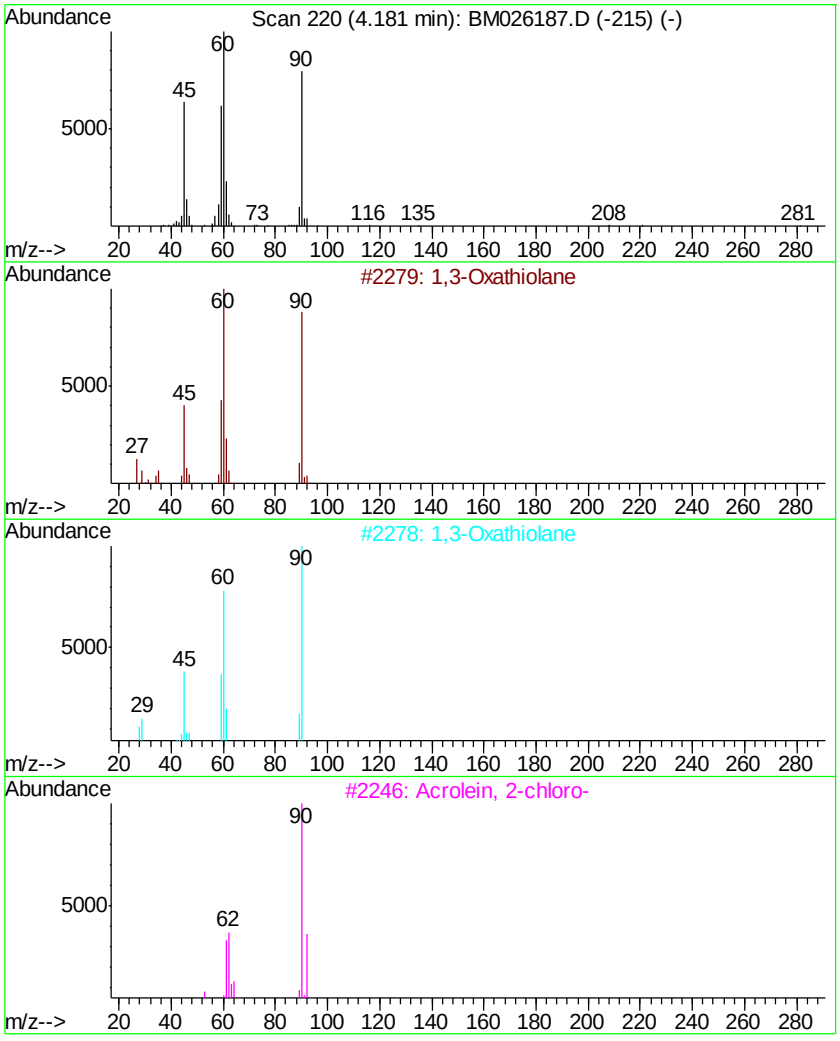
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

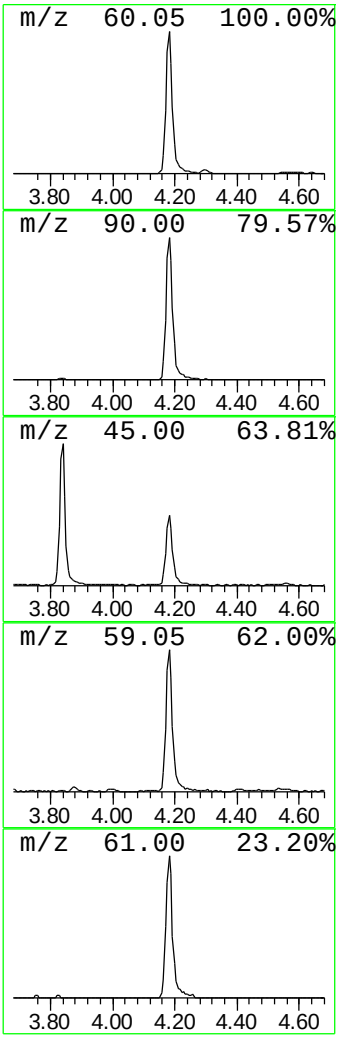
****

Peak Number 2 1,3-Oxathiolane Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.18	5.97 ng/ul	280910	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	64
3	Acrolein, 2-chloro-	90	C3H3ClO	000683-51-2	42
4	Hydrazinecarboxylic acid, methyl...	90	C2H6N2O2	006294-89-9	38
5	Hydrazinecarboxylic acid, methyl...	90	C2H6N2O2	006294-89-9	38





Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM052920\
Data File : BM026187.D
Acq On : 30 May 2020 00:46
Operator : MA/SJ
Sample : L2820-08
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : CB638

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

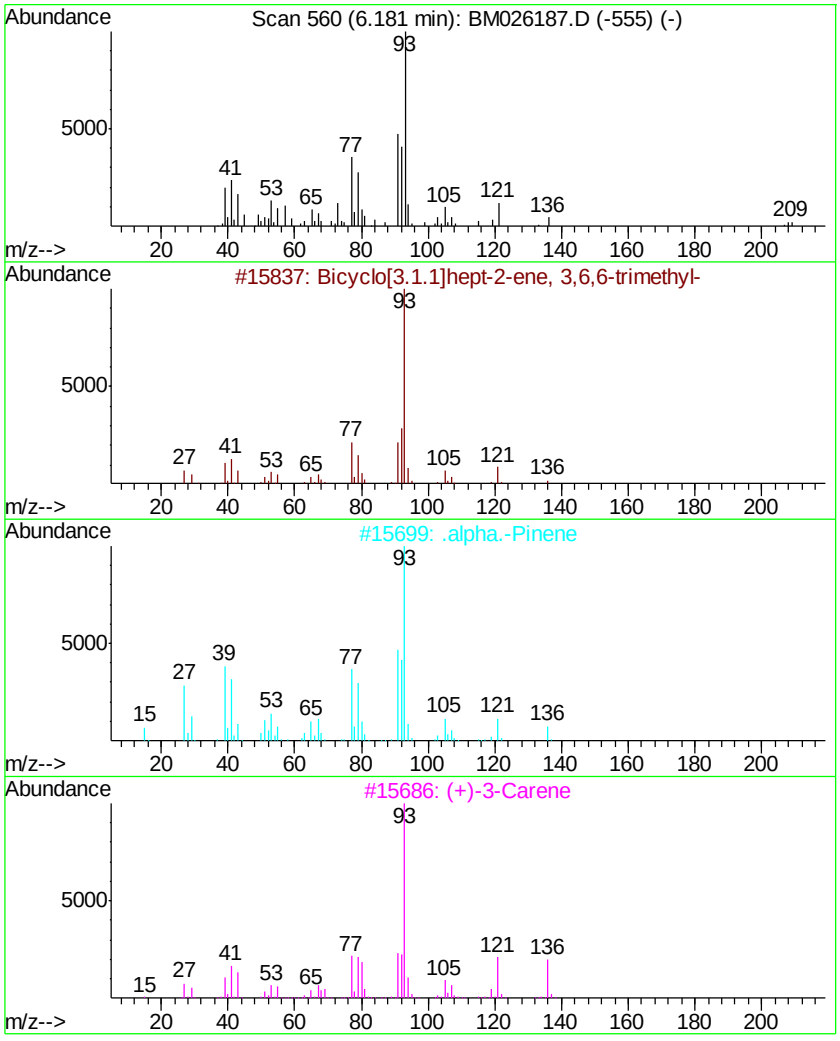
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

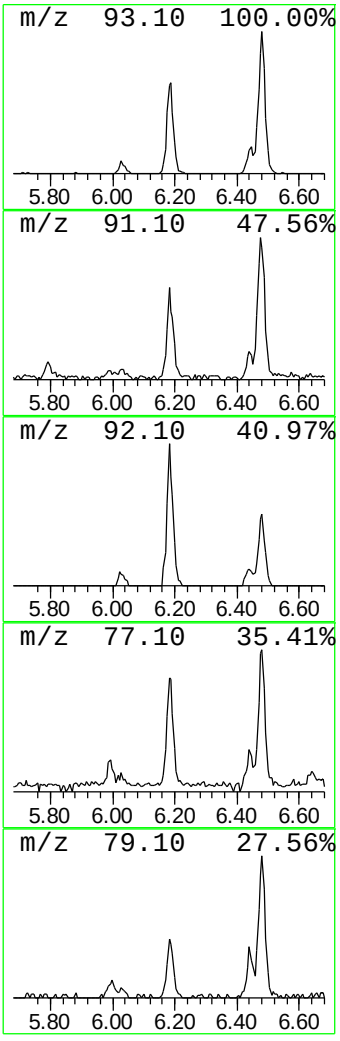
****

Peak Number 3 Bicyclo[3.1.1]hept-2-ene, 3... Concentration Rank 28

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.18	2.15 ng/ul	100999	1,4-Dichlorobenzene-d4	7.47

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Bicyclo[3.1.1]hept-2-ene, 3,6,6-...	136	C10H16	004889-83-2	93
2	.alpha.-Pinene	136	C10H16	000080-56-8	93
3	(+)-3-Carene	136	C10H16	000498-15-7	81
4	trans-.beta.-Ocimene	136	C10H16	003779-61-1	78
5	Cyclohexene, 4-methylene-1-(1-me...	136	C10H16	000099-84-3	76





Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM052920\
Data File : BM026187.D
Acq On : 30 May 2020 00:46
Operator : MA/SJ
Sample : L2820-08
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : CB638

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

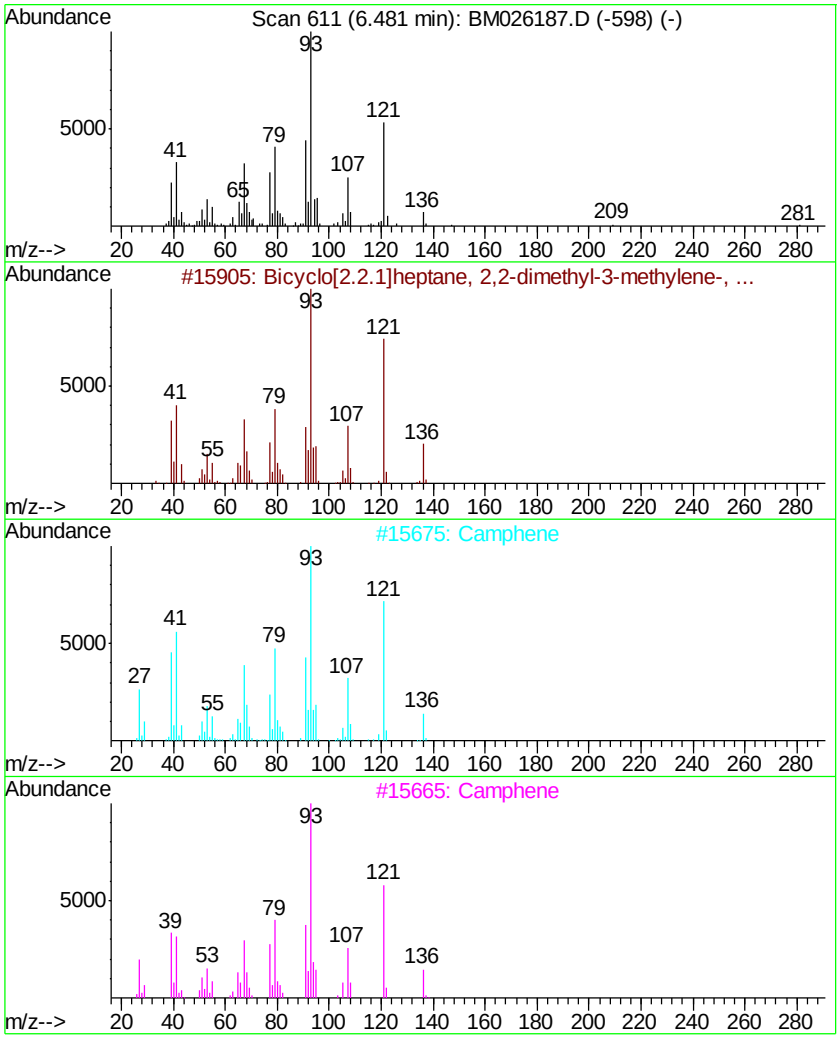
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

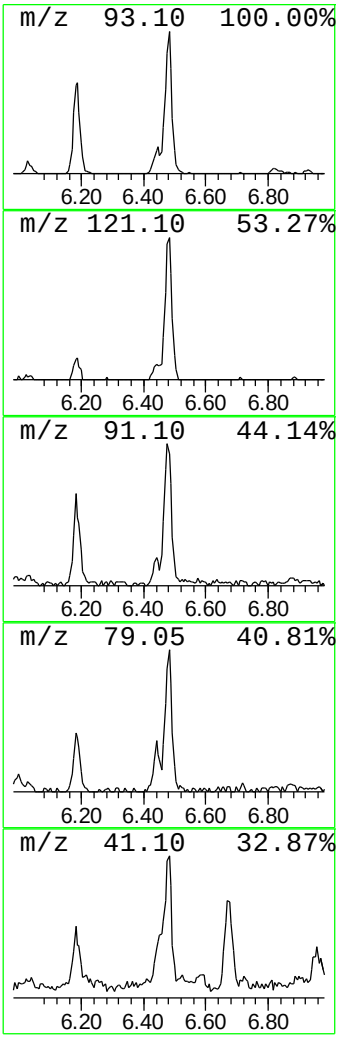
****

Peak Number 4 Bicyclo[2.2.1]heptane, 2,2-... Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.48	4.97 ng/ul	234027	1,4-Dichlorobenzene-d4	7.47

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Bicyclo[2.2.1]heptane, 2,2-dimet...	136	C10H16	005794-04-7	95
2	Camphene	136	C10H16	000079-92-5	95
3	Camphene	136	C10H16	000079-92-5	95
4	Santolina triene	136	C10H16	002153-66-4	90
5	Cyclohexene, 1-methyl-4-(1-methy...	136	C10H16	000586-62-9	74





Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM052920\
Data File : BM026187.D
Acq On : 30 May 2020 00:46
Operator : MA/SJ
Sample : L2820-08
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : CB638

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

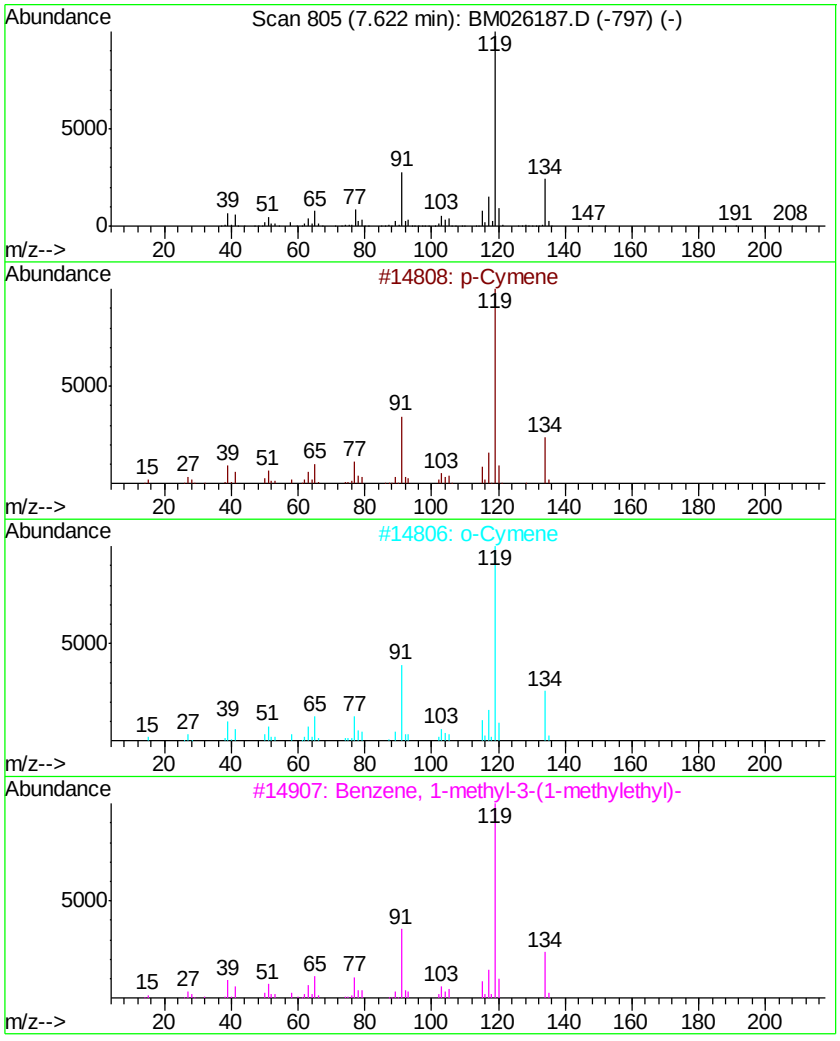
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

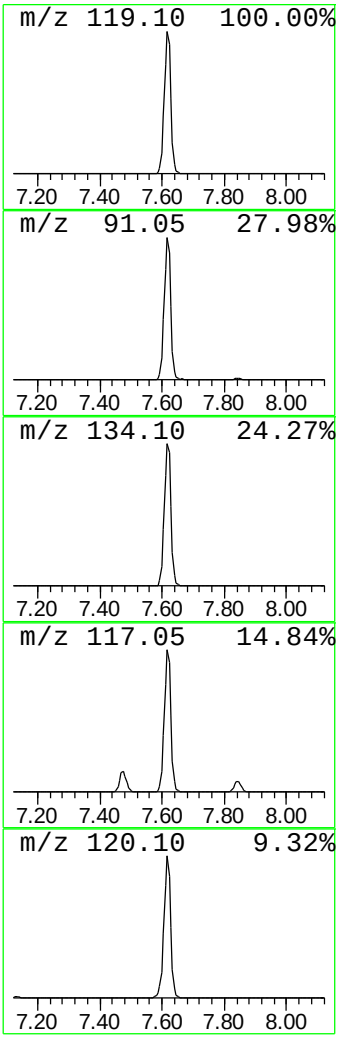
****

Peak Number 5 p-Cymene Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.62	72.21 ng/ul	3398970	1,4-Dichlorobenzene-d4	7.47

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	p-Cymene	134	C10H14	000099-87-6	97
2	o-Cymene	134	C10H14	000527-84-4	96
3	Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	95
4	o-Cymene	134	C10H14	000527-84-4	94
5	o-Cymene	134	C10H14	000527-84-4	94





Library Search Compound Report

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Acq On : 30 May 2020 00:46
Operator : MA/SJ
Sample : L2820-08
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : CB638

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Quant Title : SVOA CALIBRATION

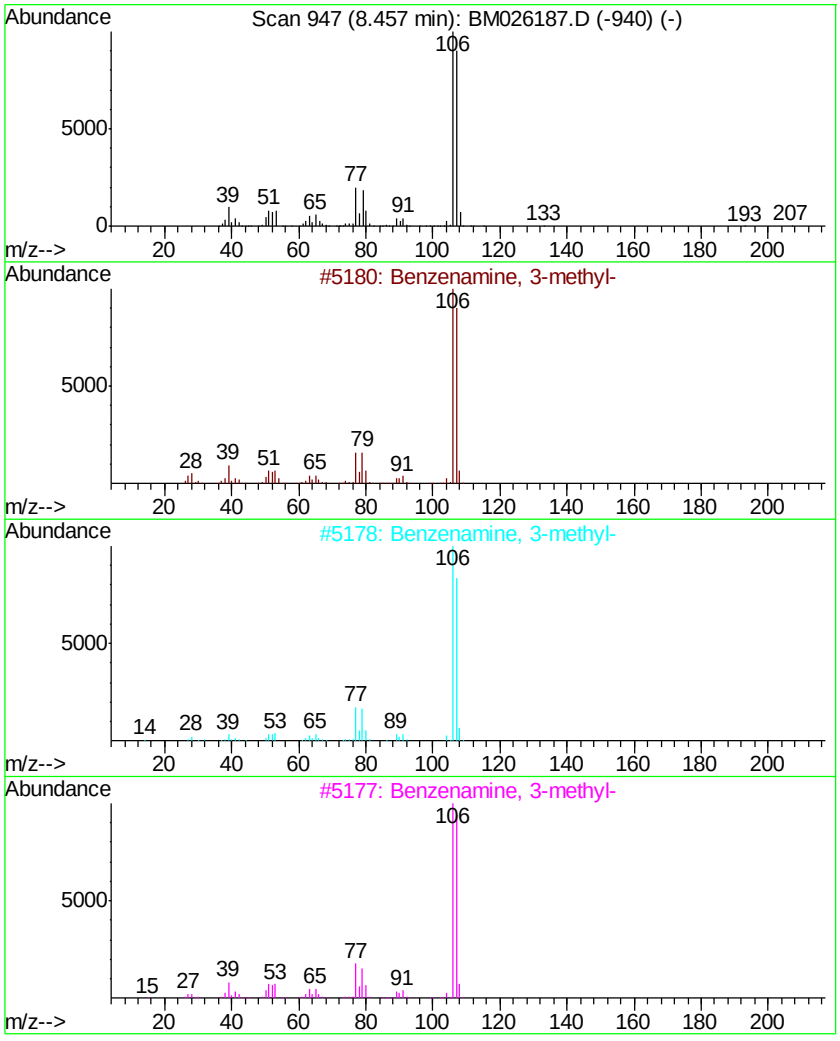
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

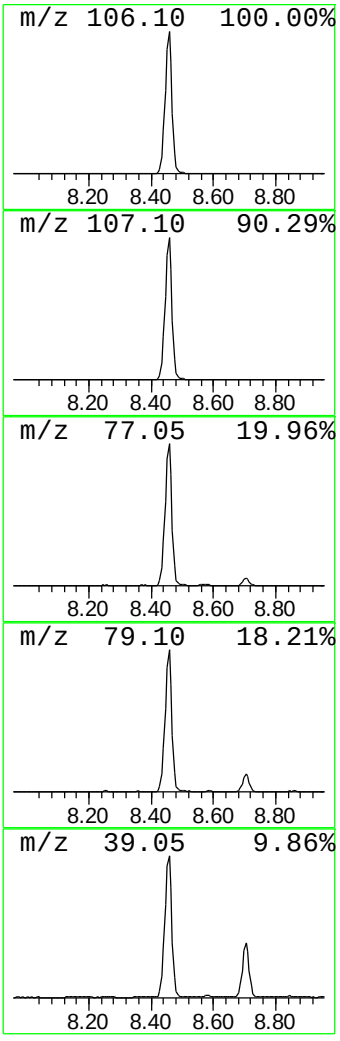
****

Peak Number 6 Benzenamine, 3-methyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.46	163.09 ng/ul	7676540	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	Benzenamine, 3-methyl-	107	C7H9N	000108-44-1	95





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Instrument : BNA_M
ClientSampleId : CB638

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Quant Title : SVOA CALIBRATION

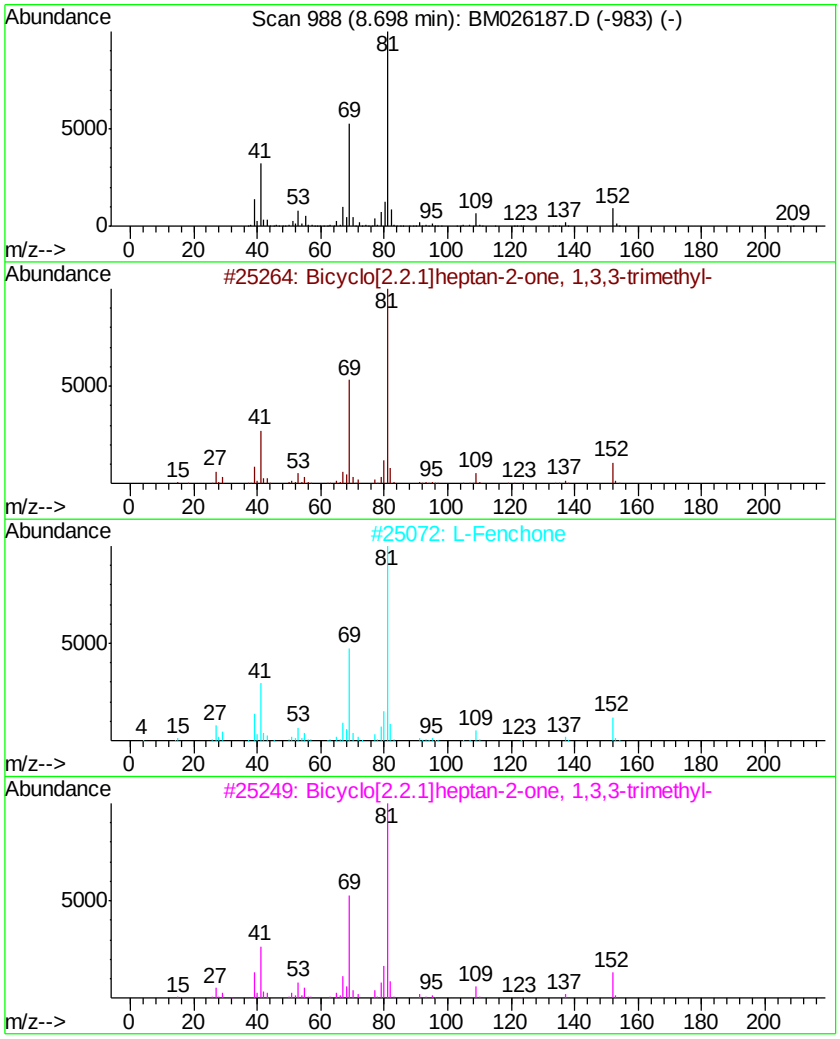
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

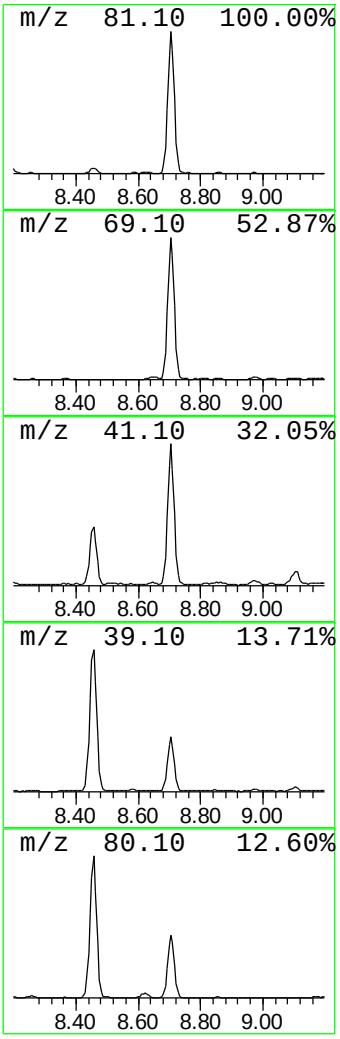
****

Peak Number 7 Bicyclo[2.2.1]heptan-2-one,... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.70	41.94 ng/ul	1974190	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	L-Fenchone	152	C10H16O	007787-20-4	95
3	Bicyclo[2.2.1]heptan-2-one, 1,3,...	152	C10H16O	001195-79-5	95
4	Bicyclo[2.2.1]heptan-2-one, 1,3,...	152	C10H16O	001195-79-5	94
5	L-Fenchone	152	C10H16O	007787-20-4	91





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Sample : L2820-08
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : CB638

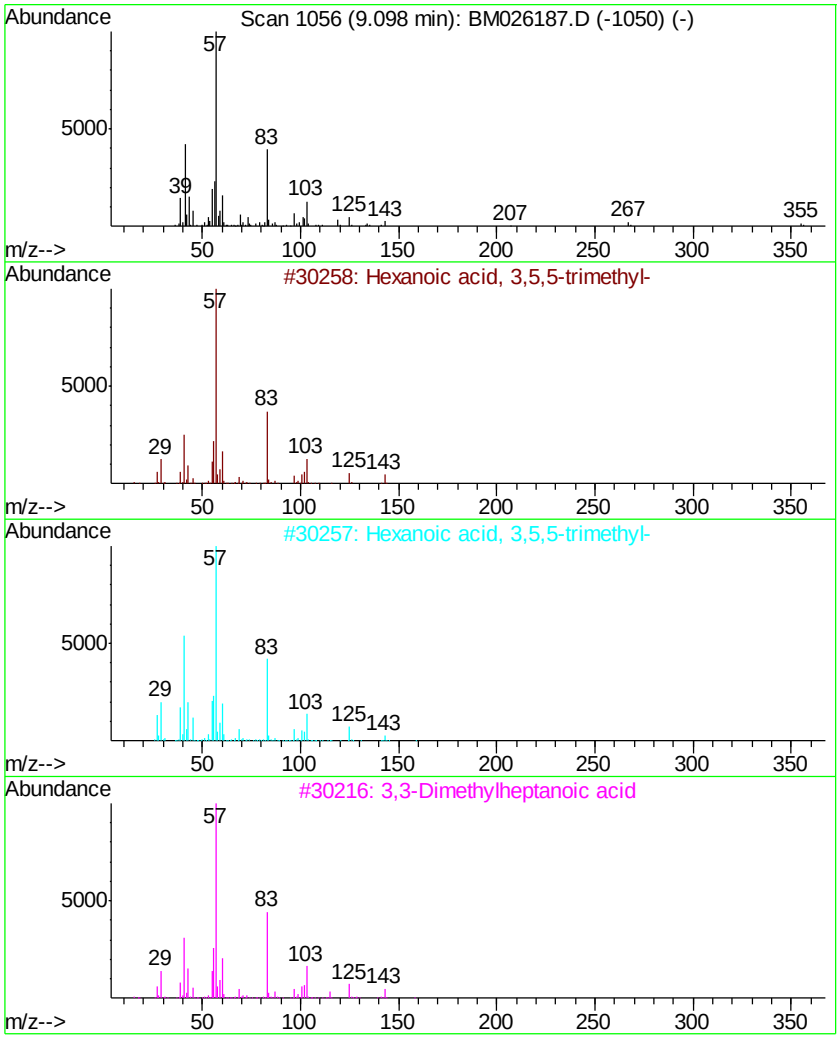
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM051320MA.M
Quant Title : SVOA CALIBRATION

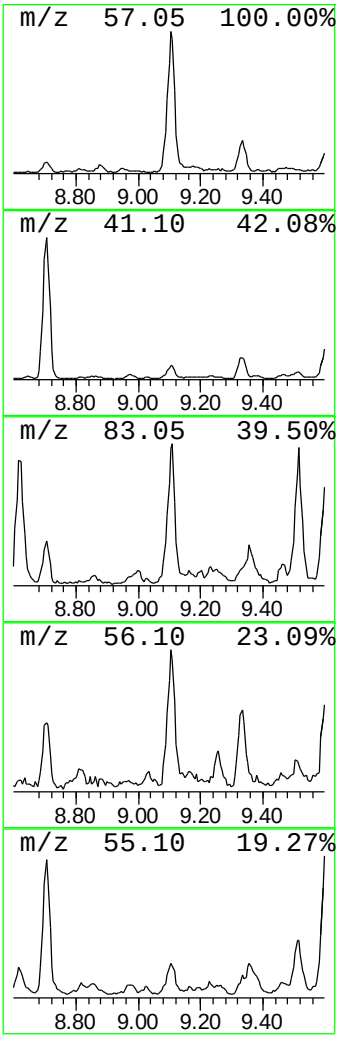
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 8 Hexanoic acid, 3,5,5-trimet... Concentration Rank 21

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.10	3.65 ng/ul	216906	Naphthalene-d8	10.23

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexanoic acid, 3,5,5-trimethyl-	158	C9H18O2	003302-10-1	90
2	Hexanoic acid, 3,5,5-trimethyl-	158	C9H18O2	003302-10-1	90
3	3,3-Dimethylheptanoic acid	158	C9H18O2	067061-30-7	90
4	Hexanoic acid, 3,5,5-trimethyl-	158	C9H18O2	003302-10-1	83
5	Hexanoyl chloride, 3,5,5-trimethyl-	176	C9H17ClO	036727-29-4	53





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Sample : L2820-08
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : CB638

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Quant Title : SVOA CALIBRATION

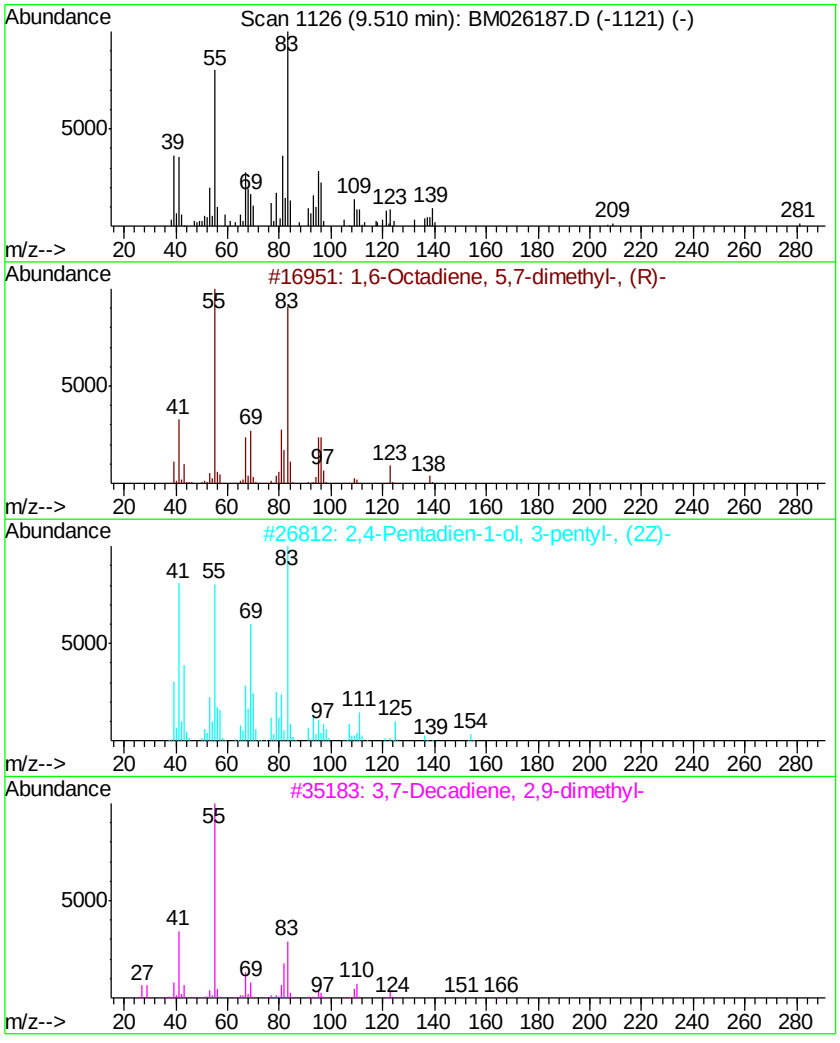
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

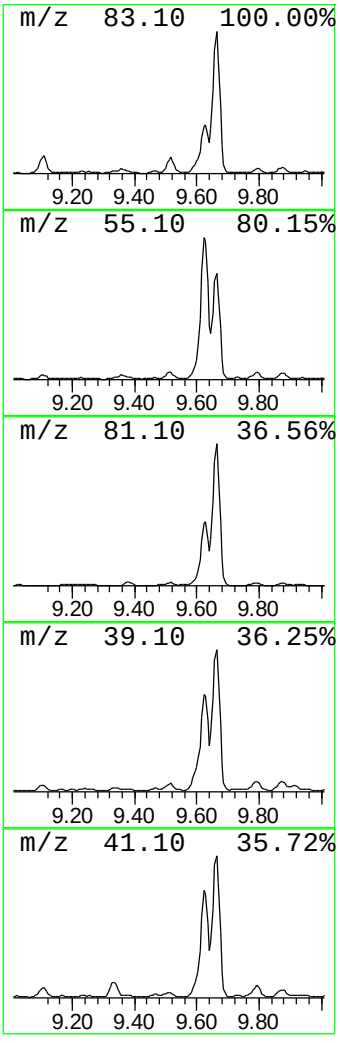
****

Peak Number 9 1,6-Octadiene, 5,7-dimethyl... Concentration Rank 23

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.51	3.15 ng/ul	187654	Naphthalene-d8	10.23

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,6-Octadiene, 5,7-dimethyl-, (R)-	138	C10H18	085006-04-8	74
2	2,4-Pentadien-1-ol, 3-pentyl-, (...	154	C10H18O	1000142-19-7	53
3	3,7-Decadiene, 2,9-dimethyl-	166	C12H22	074630-13-0	53
4	Bicyclo[3.1.1]heptan-2-one, 6,6-...	138	C9H14O	024903-95-5	52
5	1H,3H-Furo[3,4-c]furan, tetrahydro-	114	C6H10O2	005175-36-0	50





Library Search Compound Report

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Sample : L2820-08
Misc :
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Instrument : BNA_M
ClientSampleId : CB638

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Quant Title : SVOA CALIBRATION

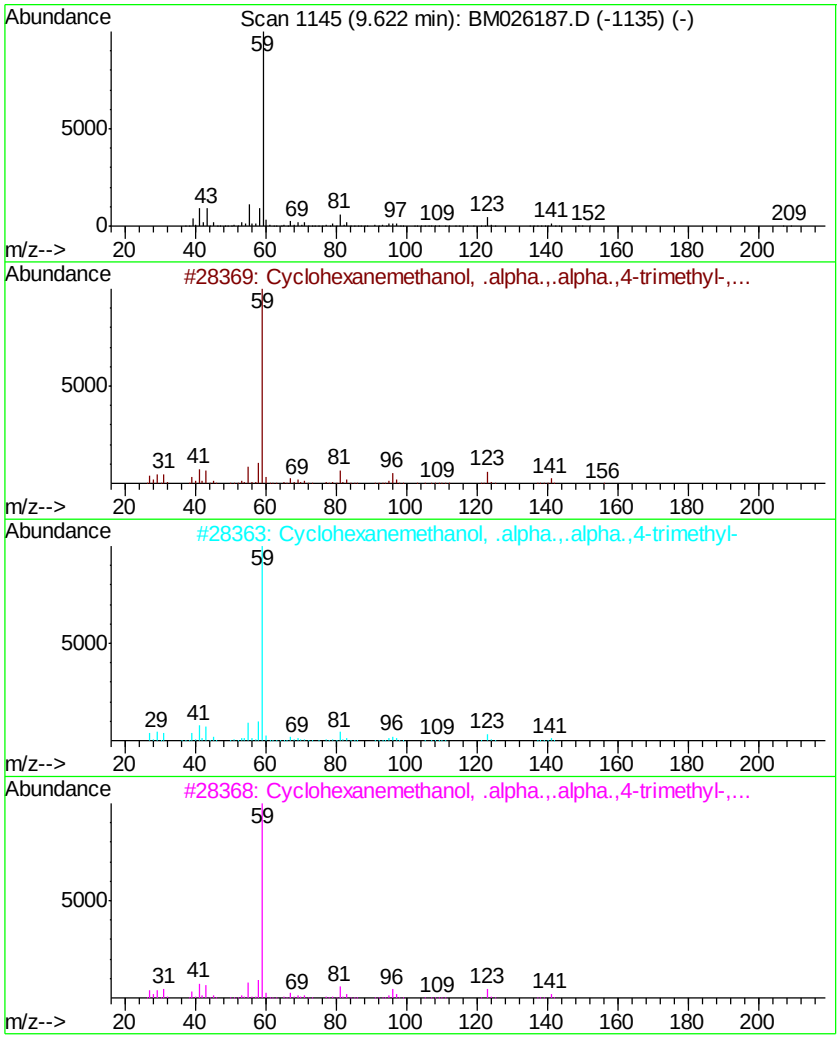
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

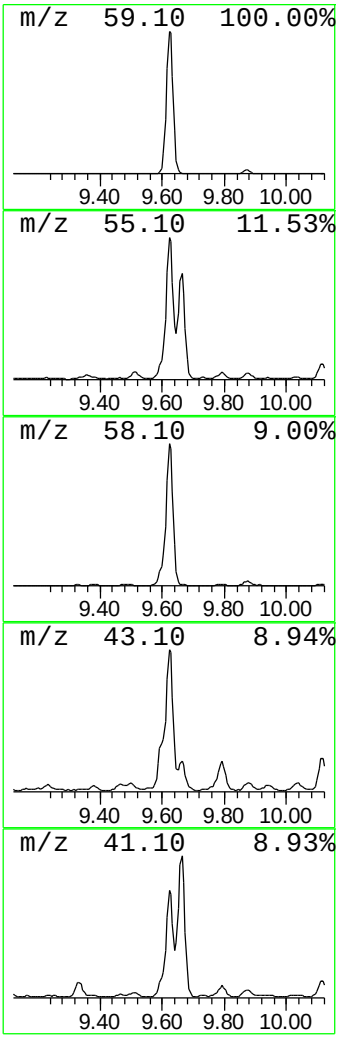
****

Peak Number 10 Cyclohexanemethanol, .alpha... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.62	96.73 ng/ul	5755920	Naphthalene-d8	10.23

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclohexanemethanol, .alpha.,.al...	156	C10H20O	005114-00-1	83
2	Cyclohexanemethanol, .alpha.,.al...	156	C10H20O	000498-81-7	83
3	Cyclohexanemethanol, .alpha.,.al...	156	C10H20O	007322-63-6	83
4	2,3-Dimethyl-4-penten-2-ol	114	C7H14O	019781-52-3	64
5	Butane, 2-methoxy-3-methyl-	102	C6H14O	062016-49-3	64





Library Search Compound Report

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Sample : L2820-08
Misc :
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Instrument : BNA_M
ClientSampleId : CB638

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Quant Title : SVOA CALIBRATION

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

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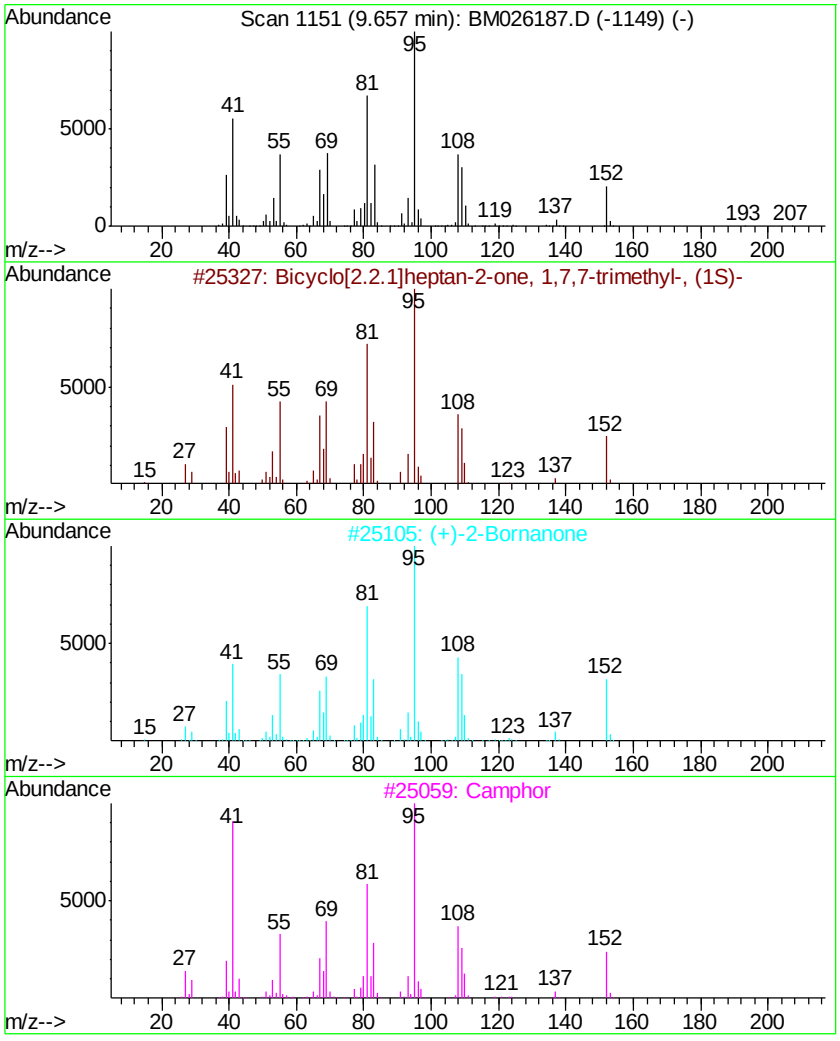
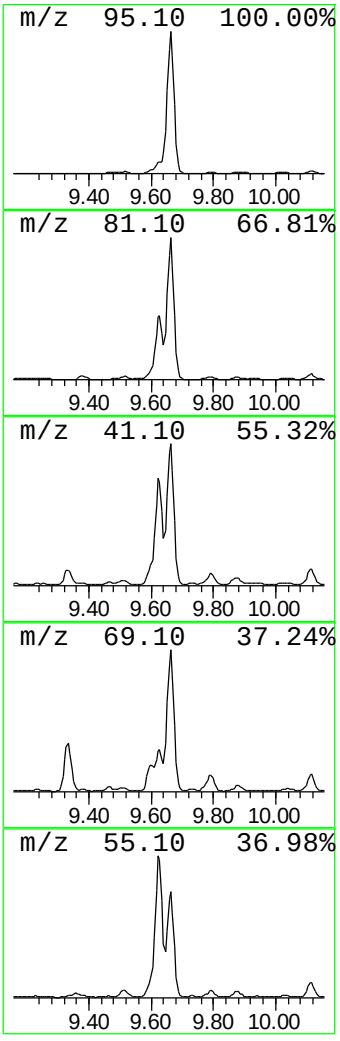
Peak Number 11 Bicyclo[2.2.1]heptan-2-one,... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.66	54.78 ng/ul	3259610	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	98
2	(+)-2-Bornanone	152	C10H16O	000464-49-3	97
3	Camphor	152	C10H16O	000076-22-2	97
4	Camphor	152	C10H16O	000076-22-2	97
5	(+)-2-Bornanone	152	C10H16O	000464-49-3	96





Library Search Compound Report

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Sample : L2820-08
Misc :
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Instrument : BNA_M
ClientSampleId : CB638

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Quant Title : SVOA CALIBRATION

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

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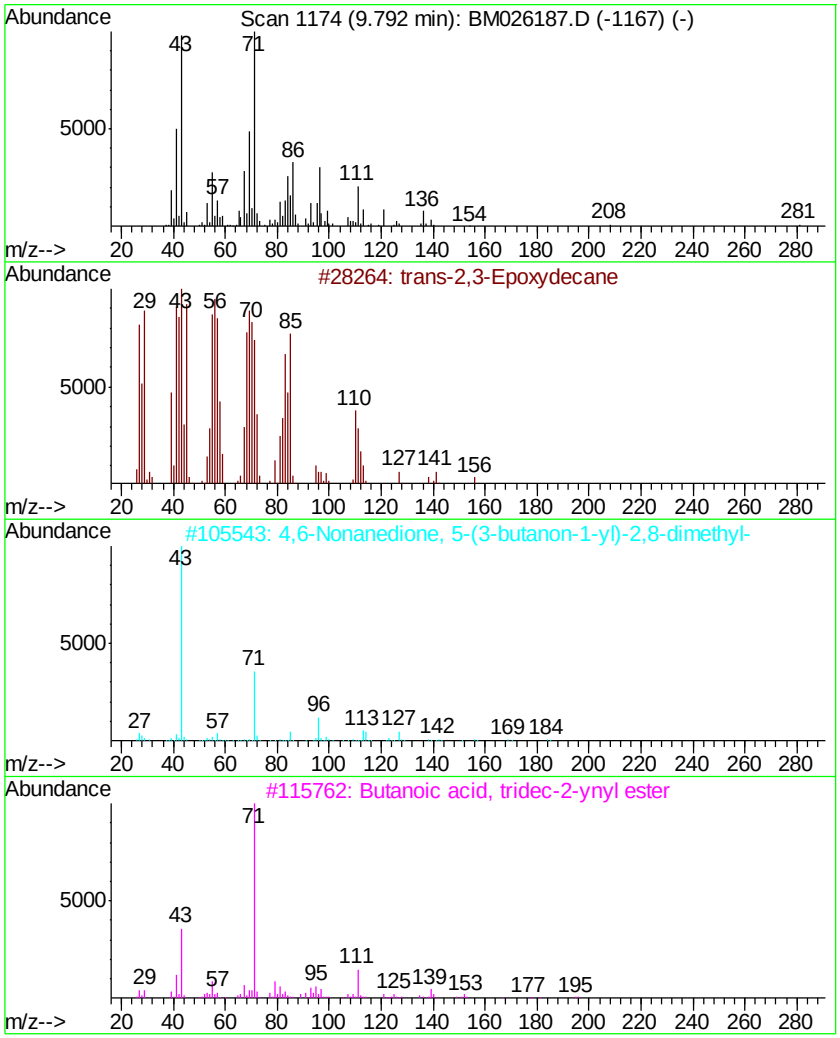
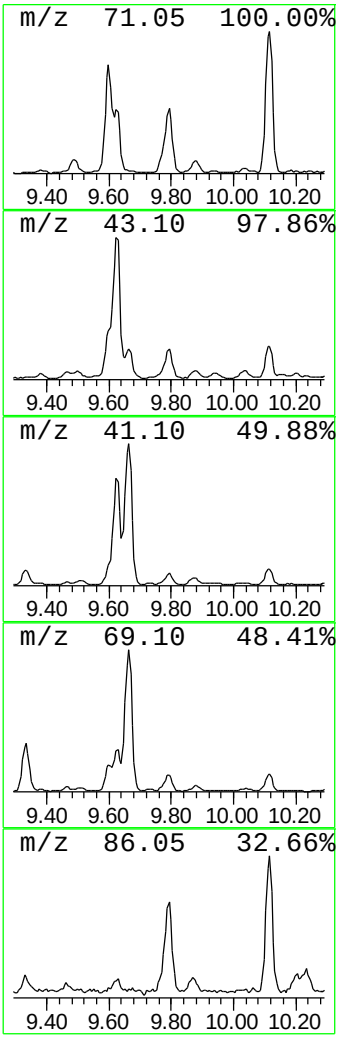
Peak Number 12 unknown-01 Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.79	6.47 ng/ul	385074	Naphthalene-d8	10.23

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	trans-2,3-Epoxydecane	156	C10H20O	054125-39-2	47
2	4,6-Nonanedione, 5-(3-butanon-1-...	254	C15H26O3	1000162-15-0	43
3	Butanoic acid, tridec-2-ynyl ester	266	C17H30O2	1000299-12-6	43
4	Diethylmalonic acid, octyl tetra...	356	C20H36O5	1000370-64-3	38
5	1-Propene, 1-methoxy-2-methyl-	86	C5H10O	017574-84-4	38





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Misc :
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Instrument : BNA_M
ClientSampleId : CB638

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Quant Title : SVOA CALIBRATION

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

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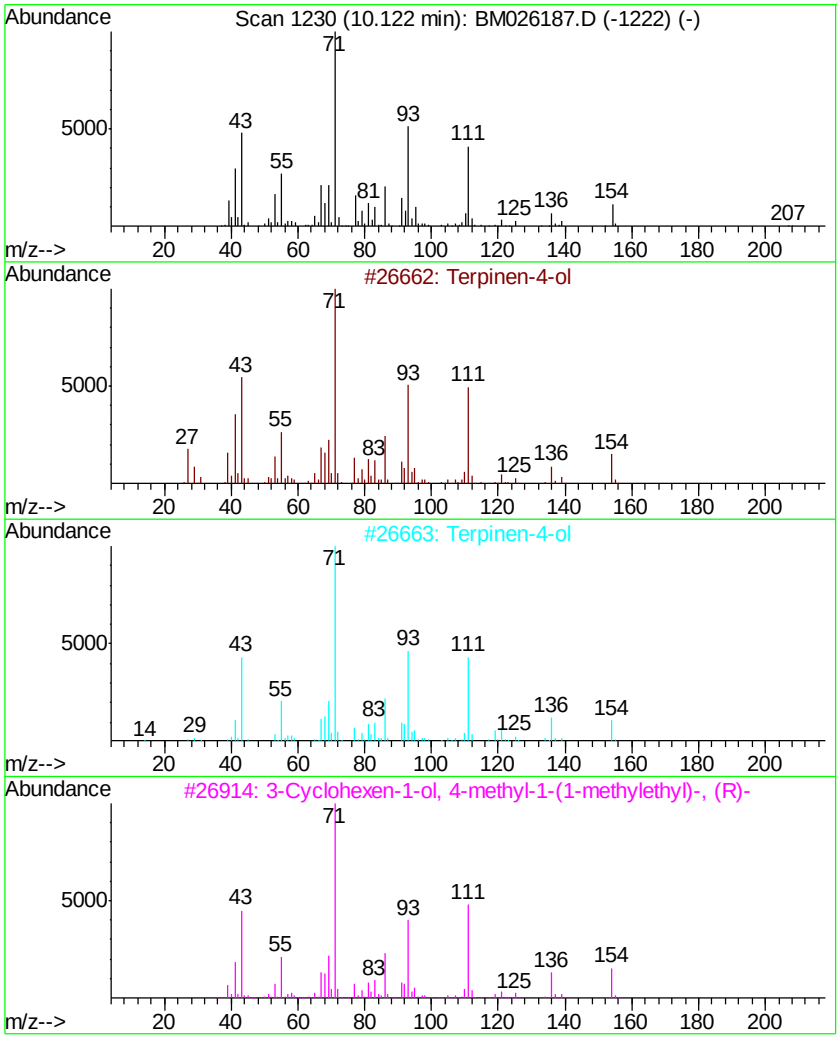
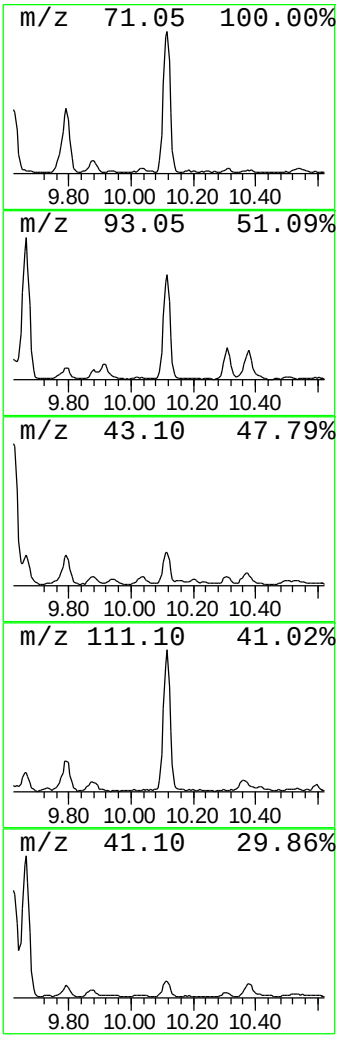
Peak Number 13 Terpinen-4-ol Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.12	10.58 ng/ul	629469	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	93
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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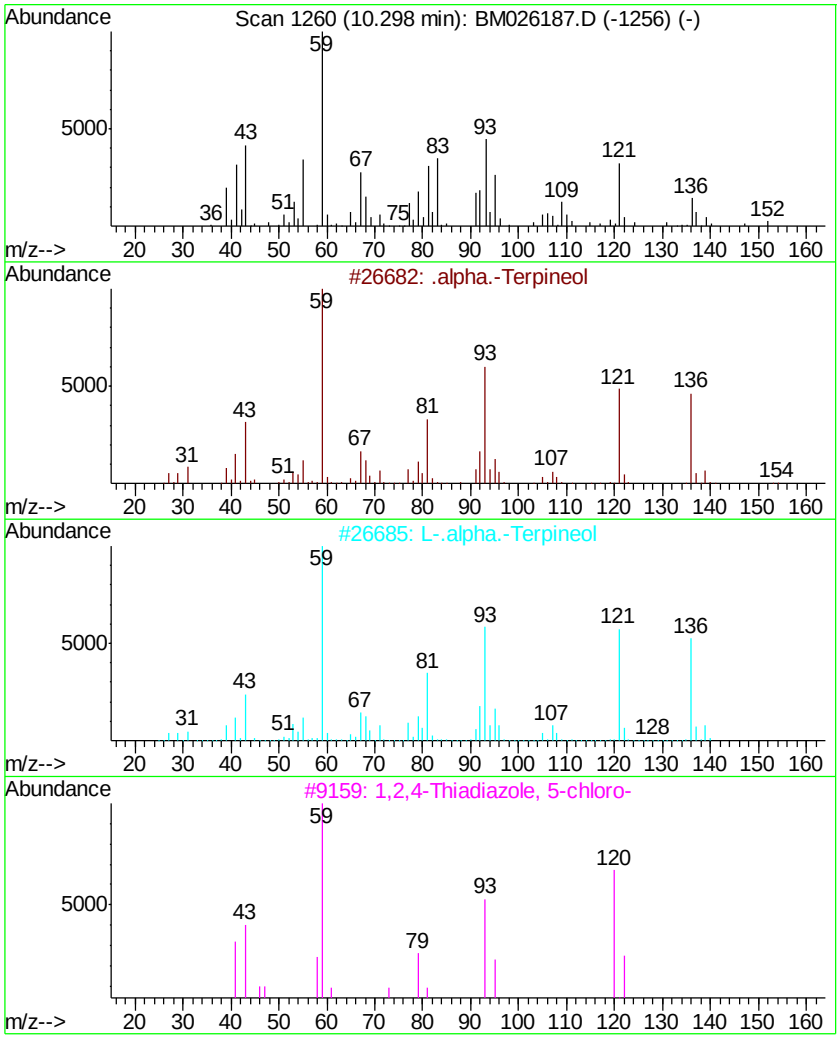
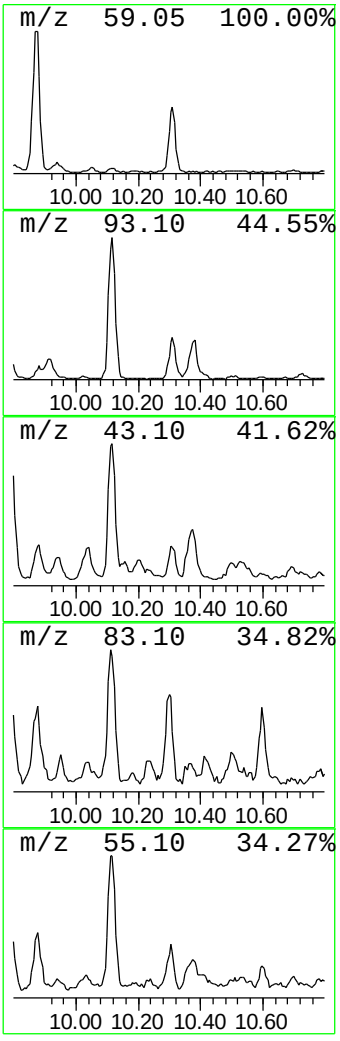
Peak Number 14 .alpha.-Terpineol Concentration Rank 22

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.30	3.35 ng/ul	199206	Naphthalene-d8	10.23

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	.alpha.-Terpineol	154	C10H18O	000098-55-5	59
2	L-.alpha.-Terpineol	154	C10H18O	010482-56-1	59
3	1,2,4-Thiadiazole, 5-chloro-	120	C2HClN2S	038362-15-1	40
4	.alpha.-Terpineol	154	C10H18O	000098-55-5	40
5	.alpha.-Terpineol	154	C10H18O	000098-55-5	38





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Sample : L2820-08
Misc :
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ClientSampleId : CB638

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Quant Title : SVOA CALIBRATION

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

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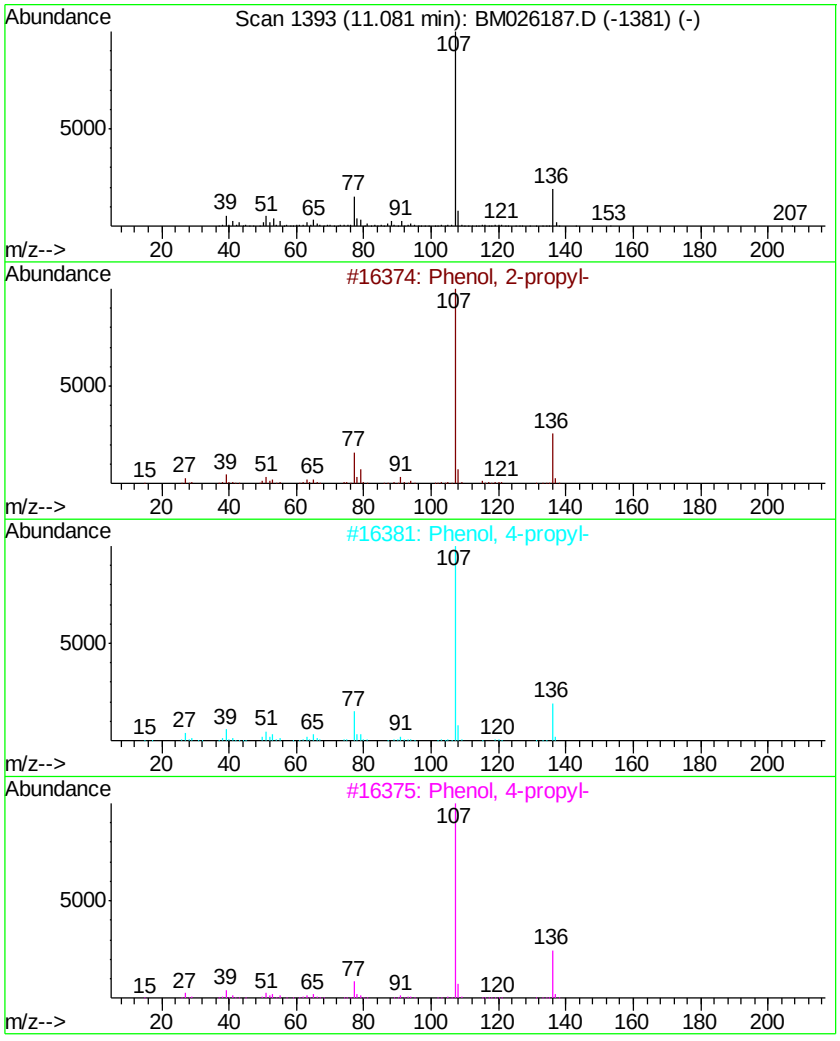
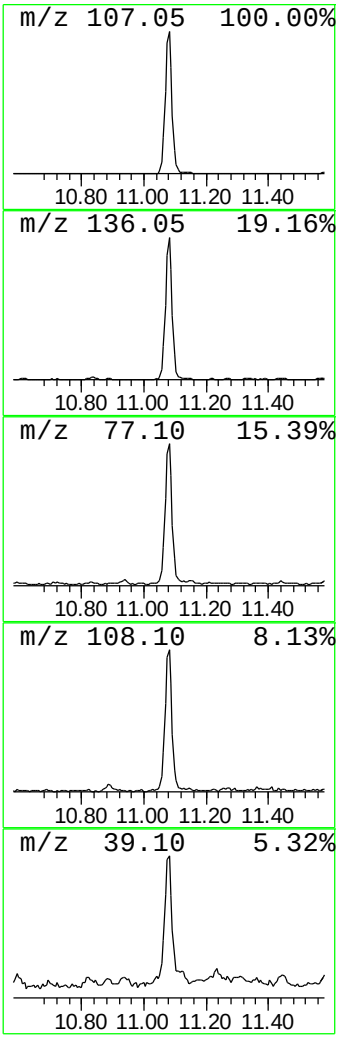
Peak Number 15 Phenol, 2-propyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.08	18.64 ng/ul	1109230	Naphthalene-d8	10.23

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, 2-propyl-	136	C9H12O	000644-35-9	91
2	Phenol, 4-propyl-	136	C9H12O	000645-56-7	91
3	Phenol, 4-propyl-	136	C9H12O	000645-56-7	87
4	Phenol, 2-propyl-	136	C9H12O	000644-35-9	78
5	Phenol, 2-propyl-	136	C9H12O	000644-35-9	72





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Instrument : BNA_M
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Quant Title : SVOA CALIBRATION

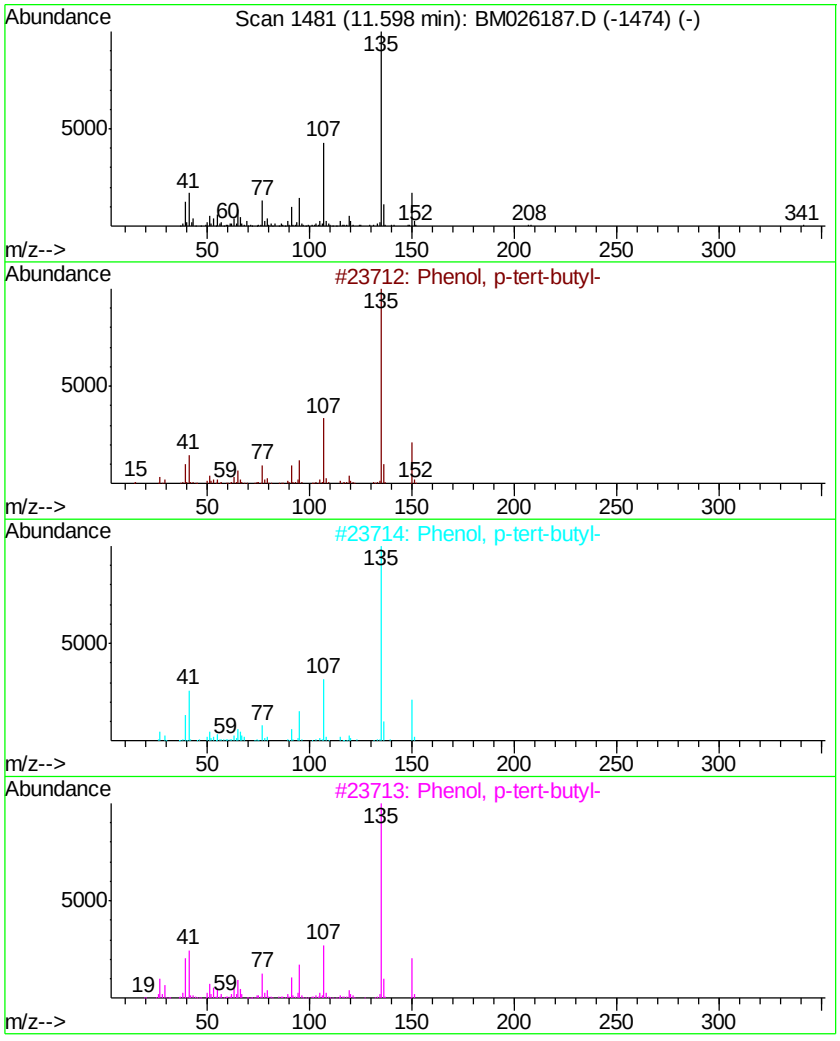
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

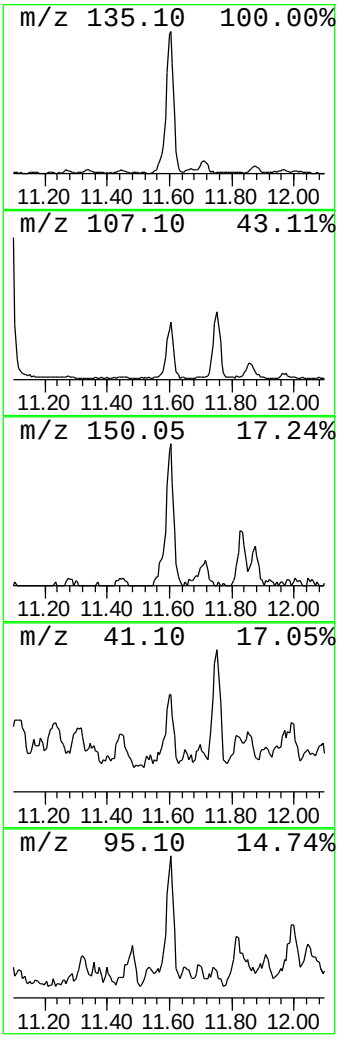
****

Peak Number 16 Phenol, p-tert-butyl- Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.60	3.82 ng/ul	227256	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, p-tert-butyl-	150	C10H14O	000098-54-4	91
3	Phenol, p-tert-butyl-	150	C10H14O	000098-54-4	80
4	3-Methoxyacetophenone	150	C9H10O2	000586-37-8	64
5	p-Anisic acid, 4-nitrophenyl ester	273	C14H11NO5	1000307-63-9	64





Library Search Compound Report

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Acq On : 30 May 2020 00:46
Operator : MA/SJ
Sample : L2820-08
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : CB638

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

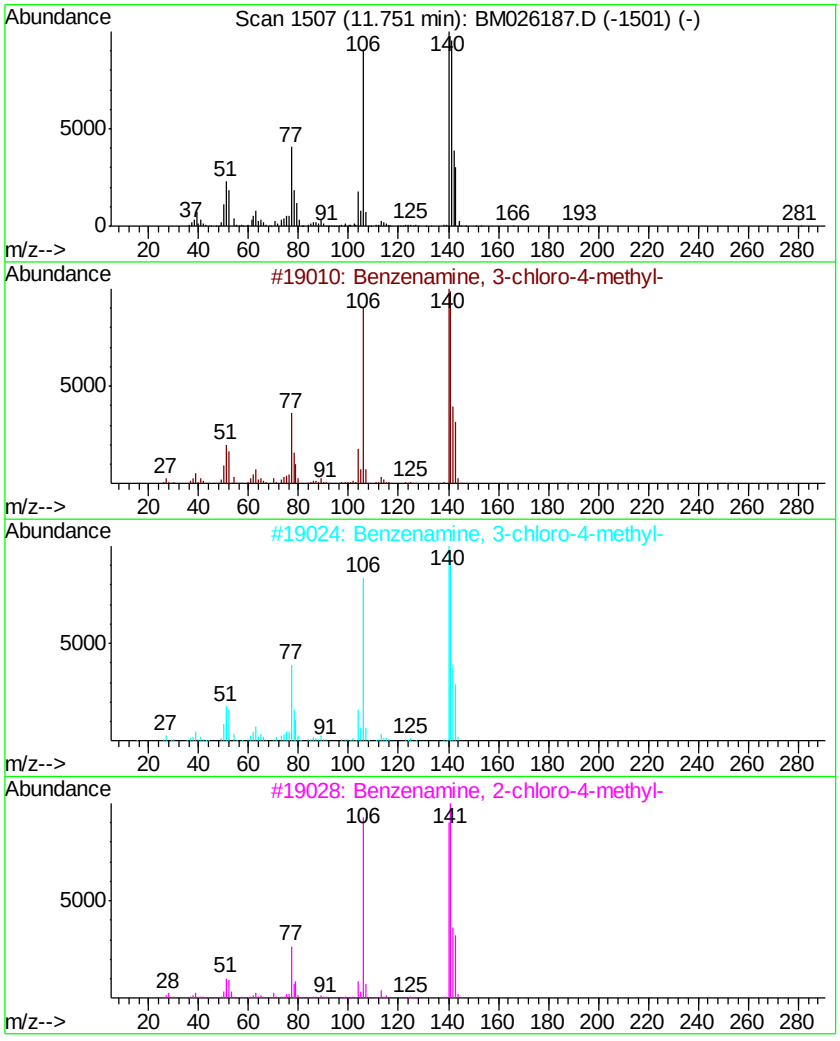
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

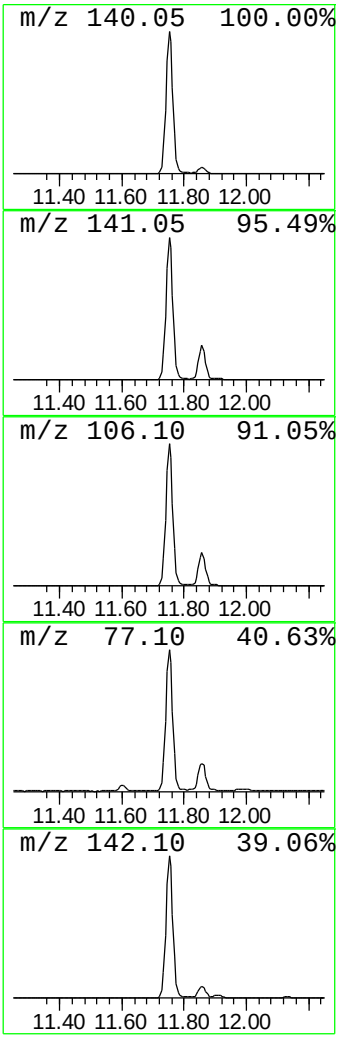
****

Peak Number 17 Benzenamine, 3-chloro-4-met... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.75	46.33 ng/ul	2756760	Naphthalene-d8	10.23

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





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Acq On : 30 May 2020 00:46
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Sample : L2820-08
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : CB638

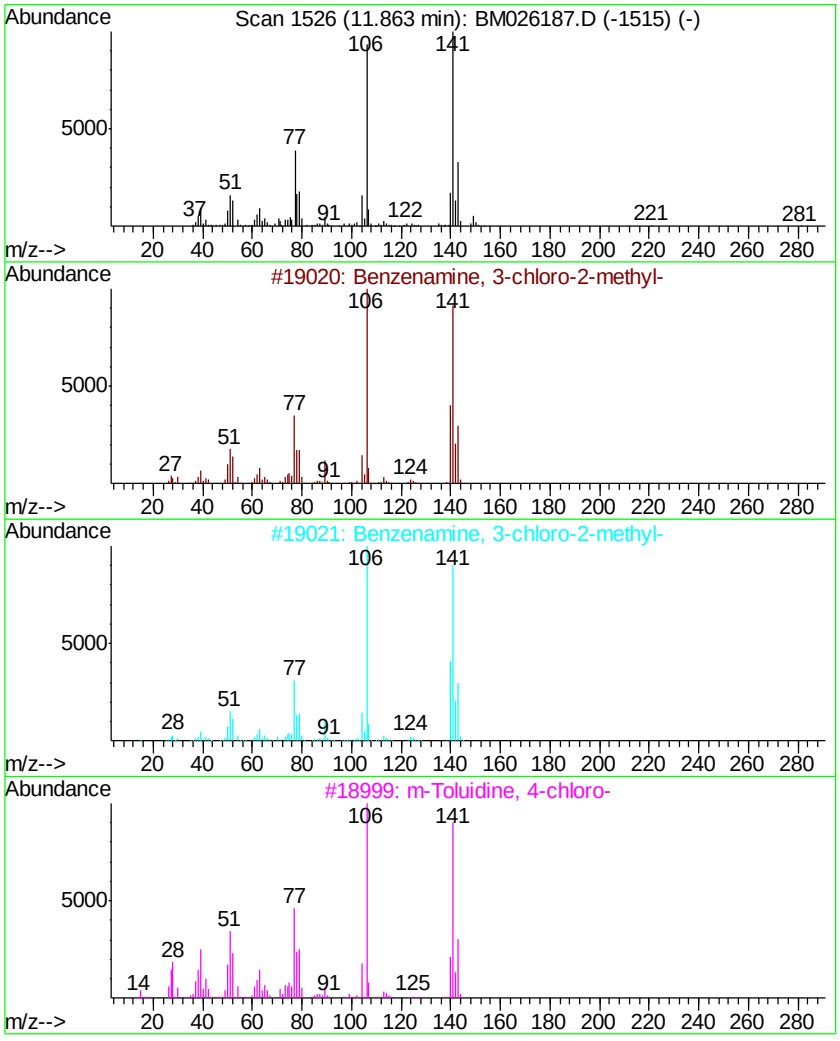
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM051320MA.M
Quant Title : SVOA CALIBRATION

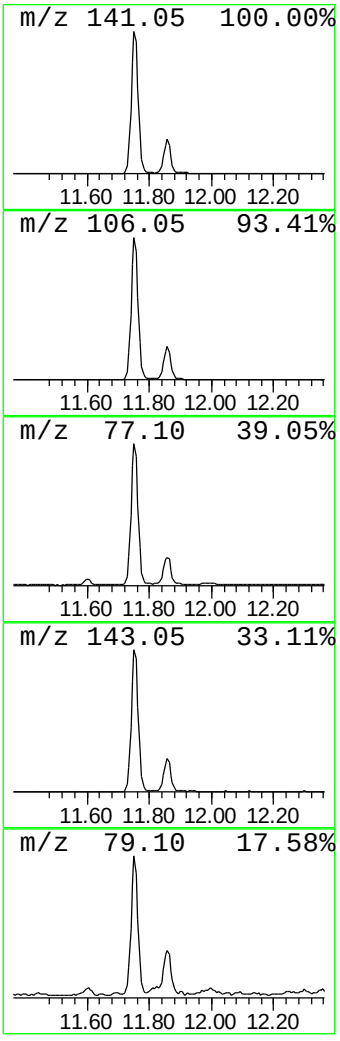
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 18 Benzenamine, 3-chloro-2-met... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.86	9.14 ng/ul	543768	Naphthalene-d8	10.23

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzenamine, 3-chloro-2-methyl-	141	C7H8ClN	000087-60-5	93
2	Benzenamine, 3-chloro-2-methyl-	141	C7H8ClN	000087-60-5	93
3	m-Toluidine, 4-chloro-	141	C7H8ClN	007149-75-9	91
4	2-Chloromethyl-5-methylpyridine	141	C7H8ClN	1000241-93-8	87
5	Benzenamine, 4-chloro-2-methyl-	141	C7H8ClN	000095-69-2	87





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Acq On : 30 May 2020 00:46
Operator : MA/SJ
Sample : L2820-08
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : CB638

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

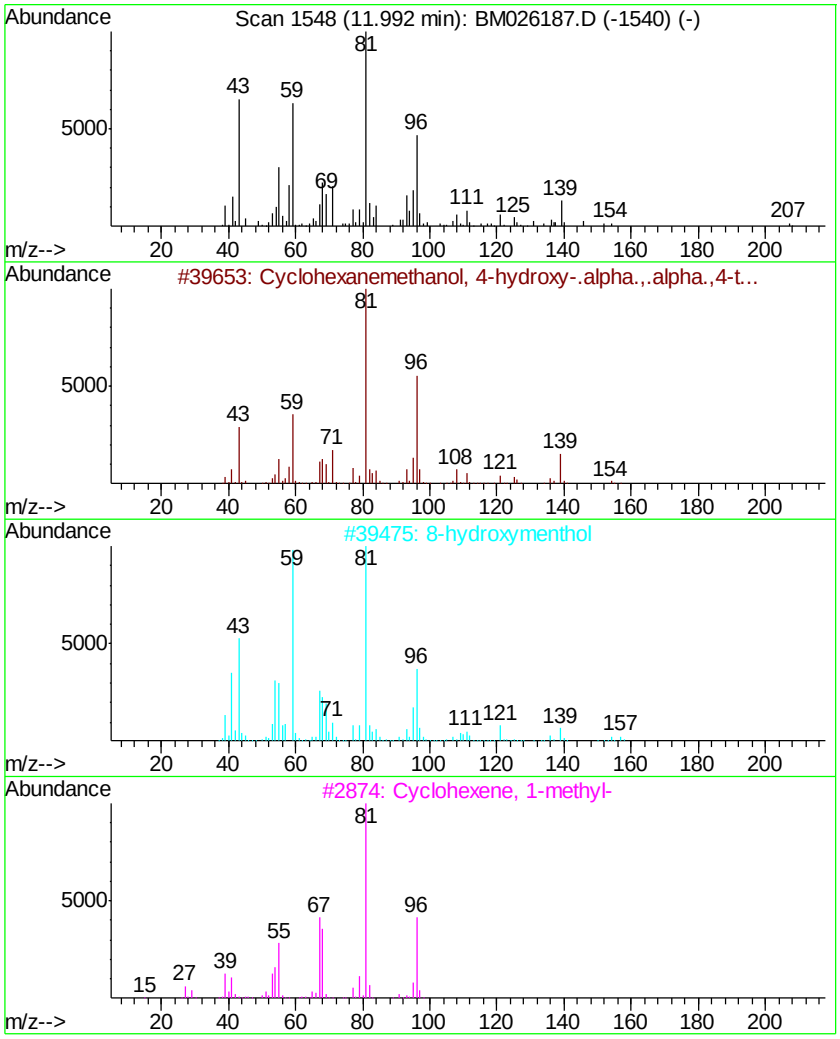
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

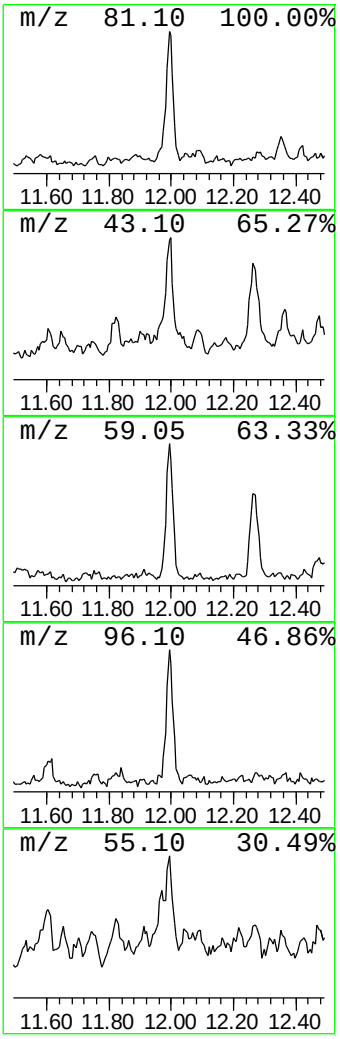
****

Peak Number 19 Cyclohexanemethanol, 4-hydr... Concentration Rank 24

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.99	3.12 ng/ul	185857	Naphthalene-d8	10.23

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclohexanemethanol, 4-hydroxy-....	172	C10H20O2	000080-53-5	64
2	8-hydroxymenthol	172	C10H20O2	1000374-16-2	50
3	Cyclohexene, 1-methyl-	96	C7H12	000591-49-1	49
4	Cyclohexane, methylene-	96	C7H12	001192-37-6	43
5	Acetic acid, trans-4-methylcyclo...	156	C9H16O2	1000368-24-6	43





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Acq On : 30 May 2020 00:46
Operator : MA/SJ
Sample : L2820-08
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : CB638

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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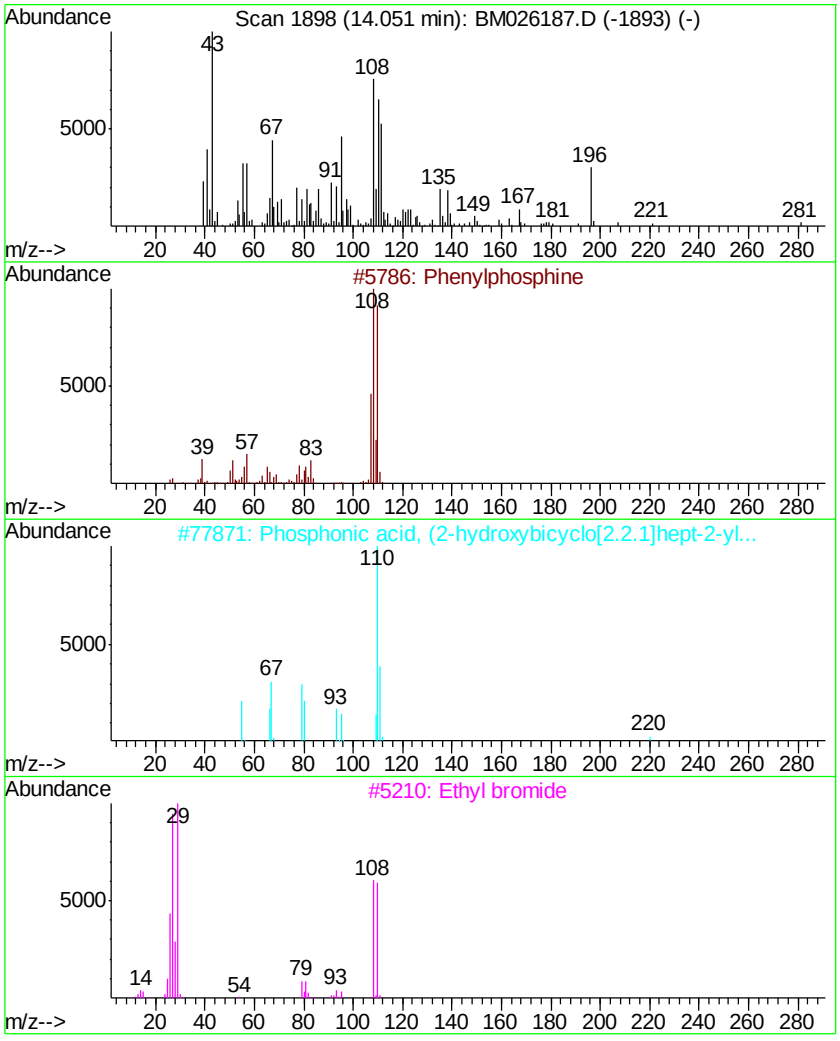
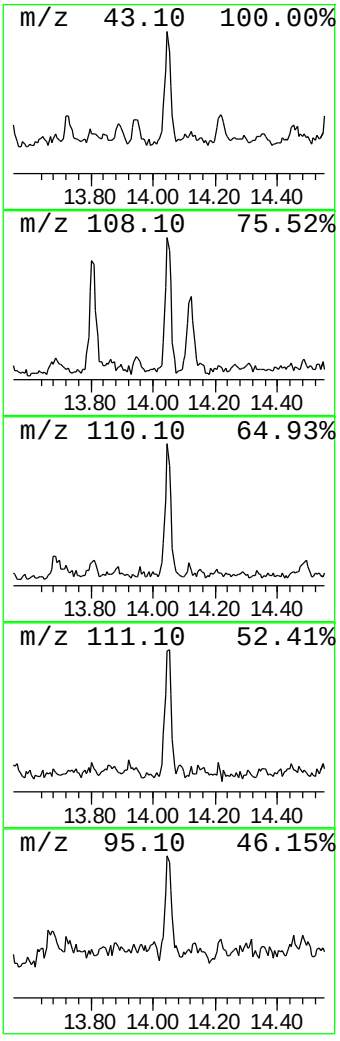
Peak Number 20 unknown-02 Concentration Rank 27

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.05	2.32 ng/ul	193817	Acenaphthene-d10	14.12

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenylphosphine	110	C6H7P	000638-21-1	27
2	Phosphonic acid, (2-hydroxybicyc...	220	C9H17O4P	027109-26-8	22
3	Ethyl bromide	108	C2H5Br	000074-96-4	22
4	Spiro[4.5]decane-1,6-dione	166	C10H14O2	036803-48-2	22
5	4-Acetonylcycloheptanone	168	C10H16O2	086428-60-6	15





Library Search Compound Report

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Data File : BM026187.D
Acq On : 30 May 2020 00:46
Operator : MA/SJ
Sample : L2820-08
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : CB638

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 1-Naphthalenol, 2-methyl- Concentration Rank 29

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.38	2.05 ng/ul	171830	Acenaphthene-d10	14.12

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Naphthalenol, 2-methyl-	158	C11H10O	007469-77-4	58
2	7-Methyl-1-naphthol	158	C11H10O	006939-33-9	52
3	1-Naphthalenemethanol	158	C11H10O	004780-79-4	49
4	1-Naphthalenemethanol	158	C11H10O	004780-79-4	46
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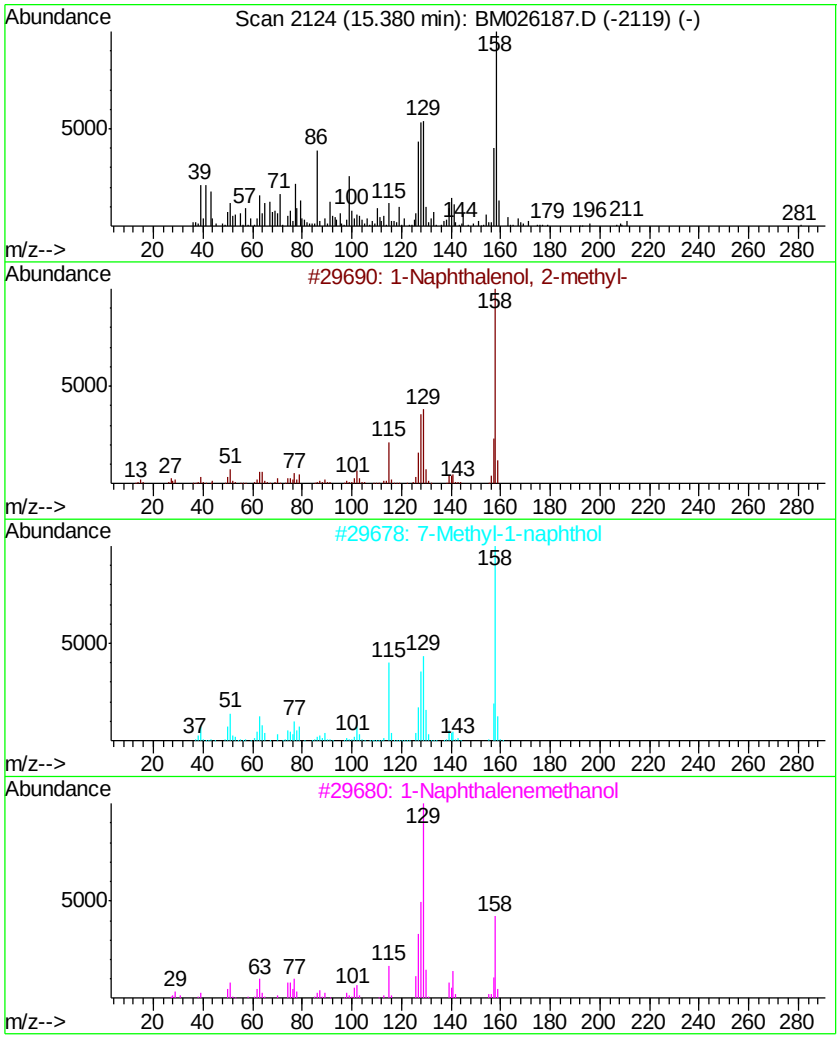

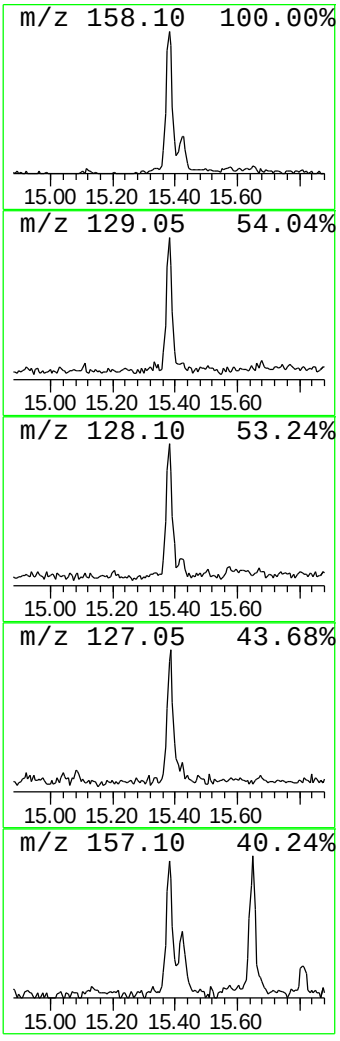

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Acq On : 30 May 2020 00:46
Operator : MA/SJ
Sample : L2820-08
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : CB638

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Quant Title : SVOA CALIBRATION

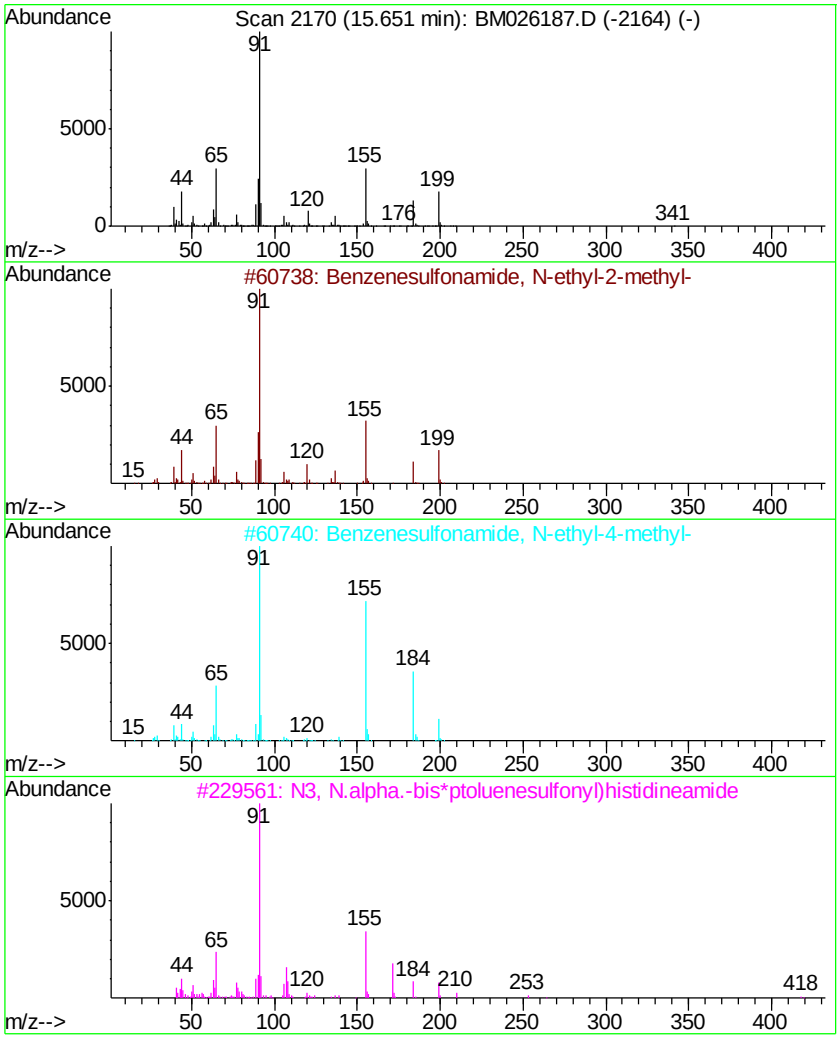
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

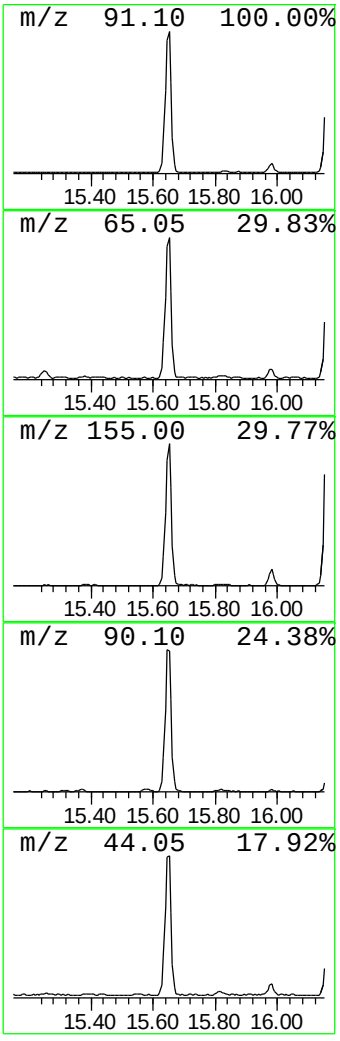
****

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.65	13.96 ng/ul	1478910	Phenanthrene-d10	16.87

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzenesulfonamide, N-ethyl-2-me...	199	C9H13NO2S	001077-56-1	98
2	Benzenesulfonamide, N-ethyl-4-me...	199	C9H13NO2S	000080-39-7	59
3	N3, N.alpha.-bis*ptoluenesulfony...	462	C20H22N4O5S2	1000130-21-9	53
4	Tolbutamide	270	C12H18N2O3S	000064-77-7	50
5	p-Toluenesulfonylacetonitrile	195	C9H9NO2S	005697-44-9	50





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Acq On : 30 May 2020 00:46
Operator : MA/SJ
Sample : L2820-08
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : CB638

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Quant Title : SVOA CALIBRATION

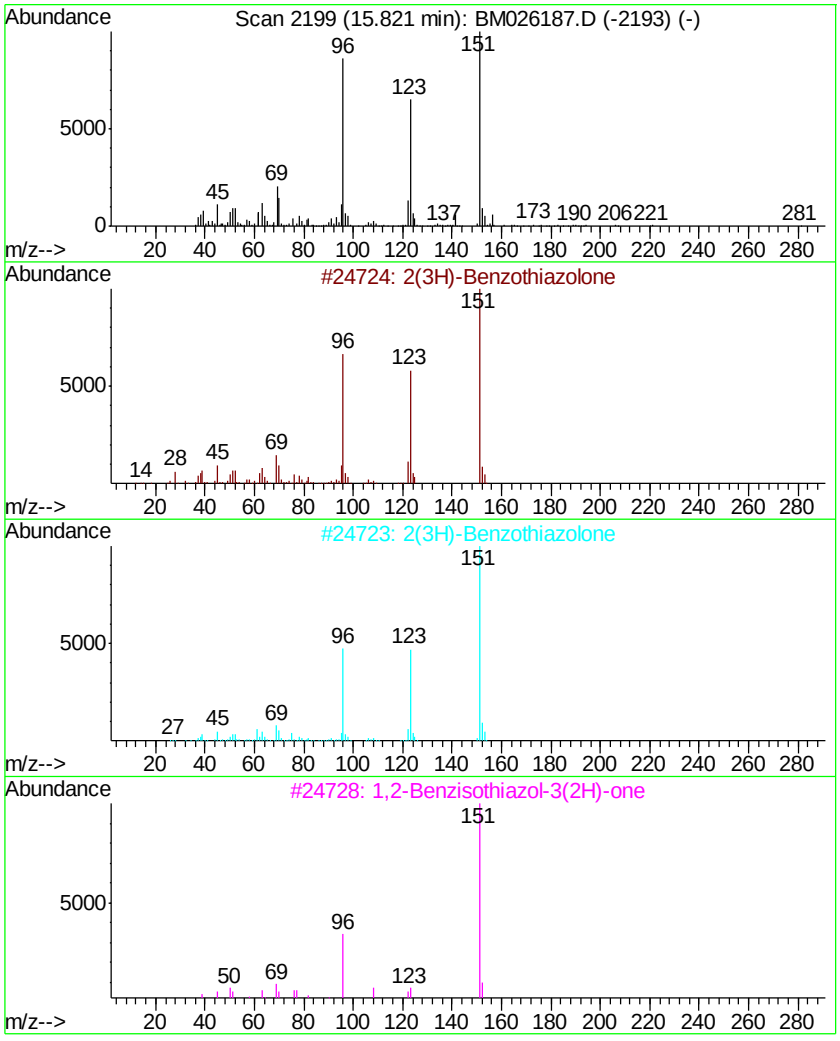
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

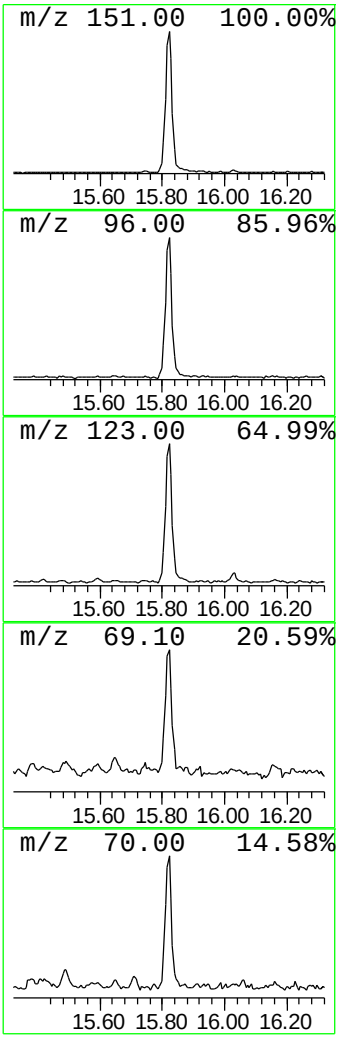
****

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.82	5.19 ng/ul	549868	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	76
3	1,2-Benzisothiazol-3(2H)-one	151	C7H5NOS	002634-33-5	76
4	4-Methoxyformanilide	151	C8H9NO2	005470-34-8	38
5	Benzooxazole-2-thiol	151	C7H5NOS	1000318-31-6	35





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Sample : L2820-08
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : CB638

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

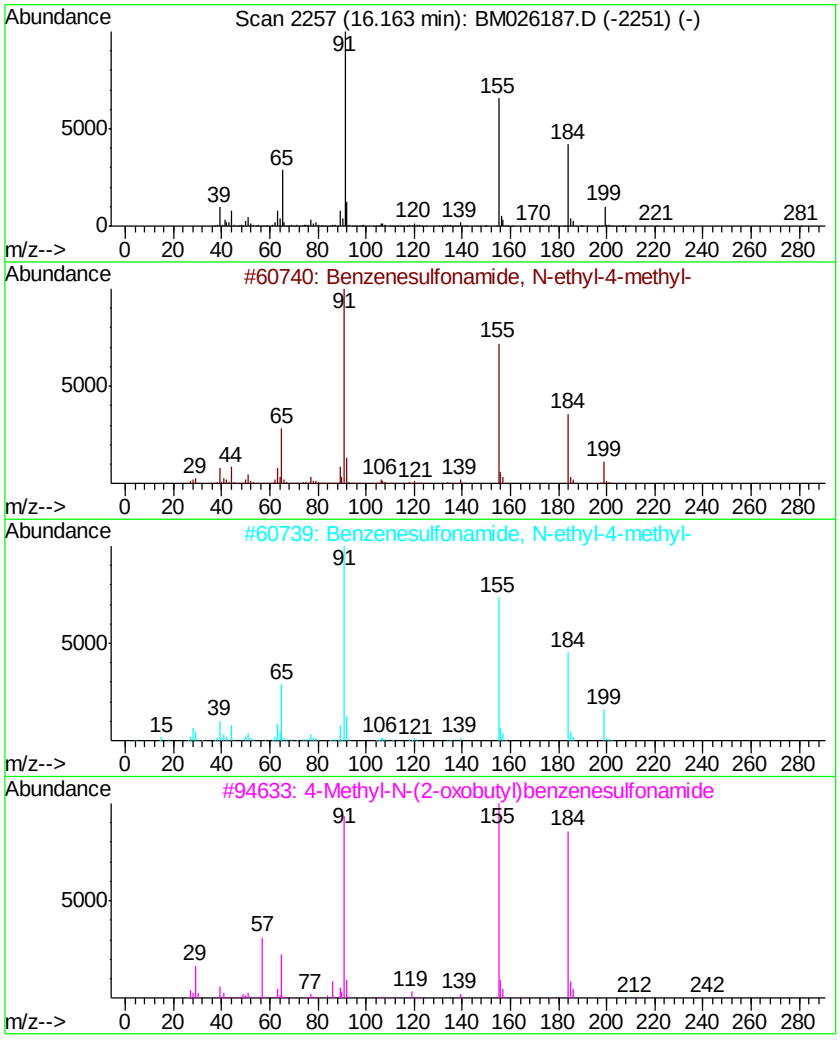
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

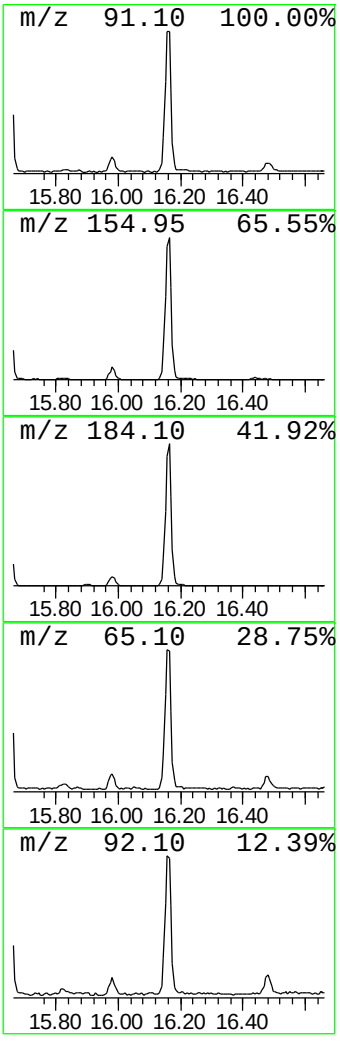
****

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.16	8.88 ng/ul	941565	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	4-Methyl-N-(2-oxobutyl)benzenesu...	241	C11H15NO3S	1000192-14-5	78
4	Furazan-3-carboxylic acid, 4-ami...	325	C12H15N5O4S	1000304-69-4	78
5	(Toluene-4-sulfonylamino)-acetic...	283	C13H17NO4S	1000190-48-0	64





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Sample : L2820-08
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : CB638

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

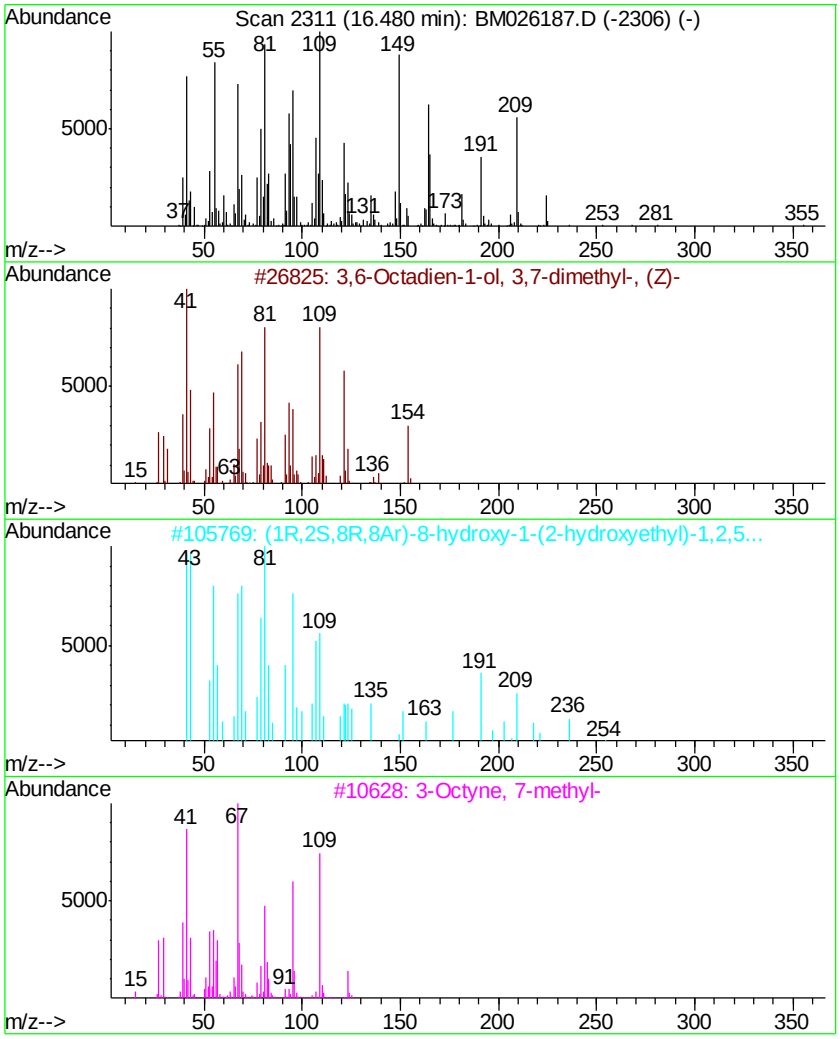
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

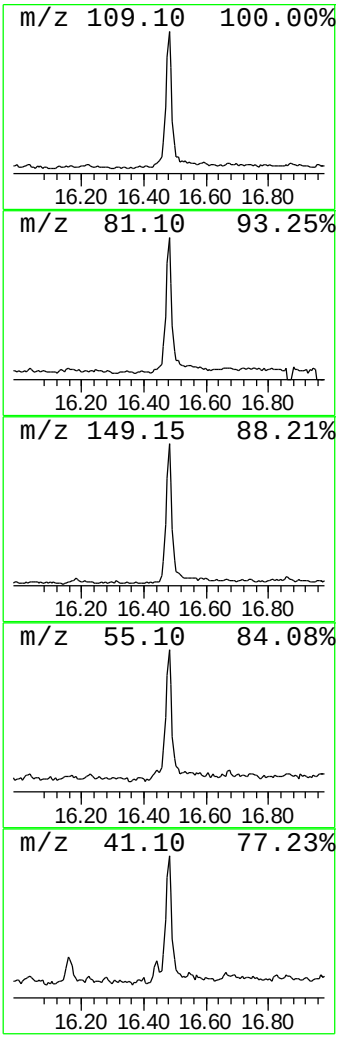
****

Peak Number 25 unknown-03 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.48	9.40 ng/ul	995673	Phenanthrene-d10	16.87

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3,6-Octadien-1-ol, 3,7-dimethyl-...	154	C10H18O	005944-20-7	46
2	(1R,2S,8R,8Ar)-8-hydroxy-1-(2-hy...	254	C16H30O2	1000298-98-3	43
3	3-Octyne, 7-methyl-	124	C9H16	037050-06-9	30
4	Cyclohexane, ethylidene-	110	C8H14	001003-64-1	18
5	Cyclohexene,3-(1-methylpropyl)-	138	C10H18	015232-91-4	14





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Sample : L2820-08
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : CB638

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

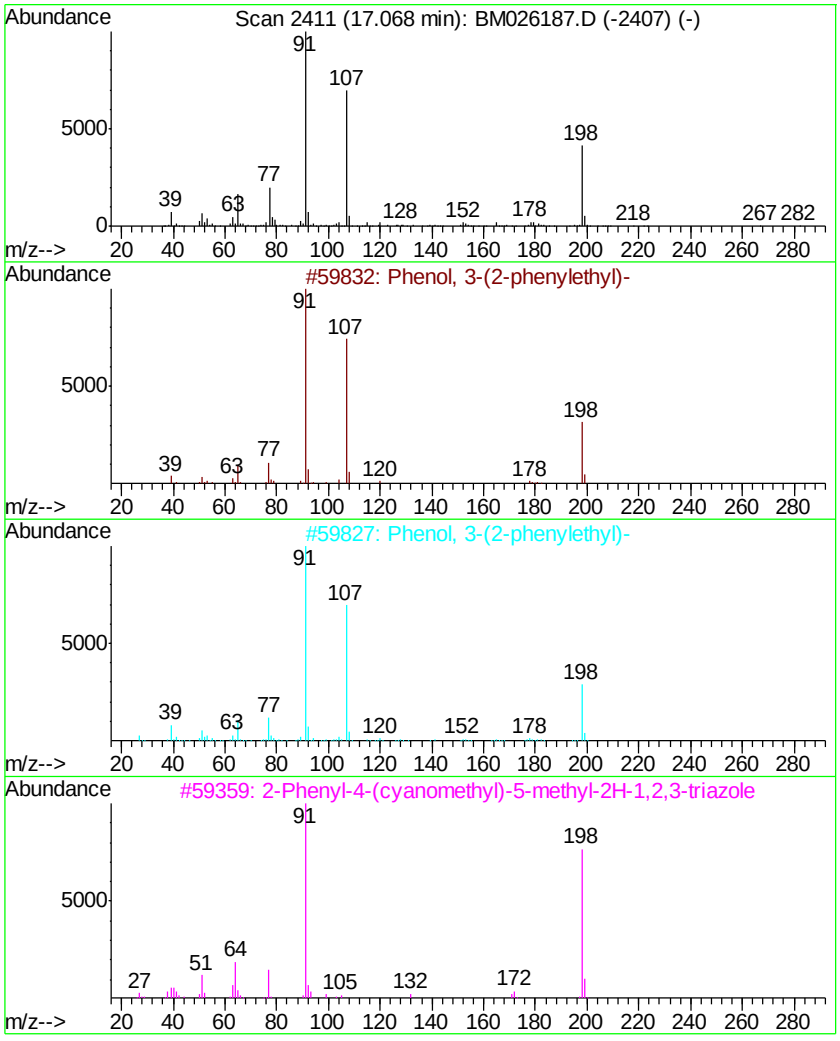
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

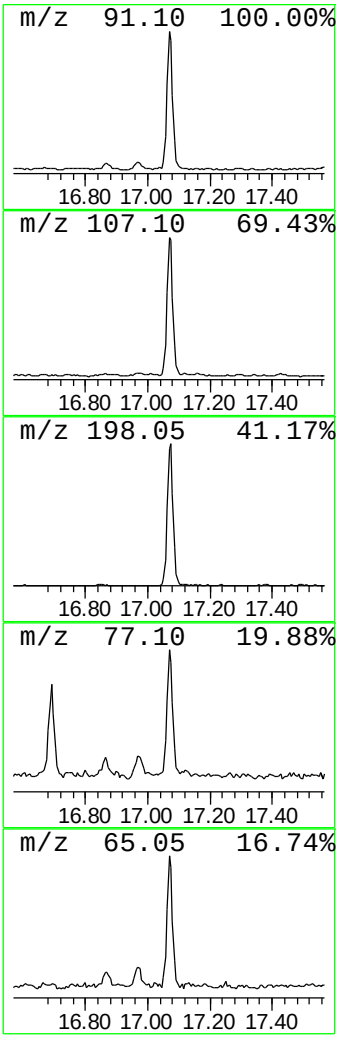
****

Peak Number 26 Phenol, 3-(2-phenylethyl)- Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.07	4.26 ng/ul	451034	Phenanthrene-d10	16.87

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	2-Phenyl-4-(cyanomethyl)-5-methy...	198	C11H10N4	013322-20-8	38
4	Pyridine, 2,3-dimethyl-	107	C7H9N	000583-61-9	30
5	Pyridine, 2,3-dimethyl-	107	C7H9N	000583-61-9	27





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Operator : MA/SJ
Sample : L2820-08
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : CB638

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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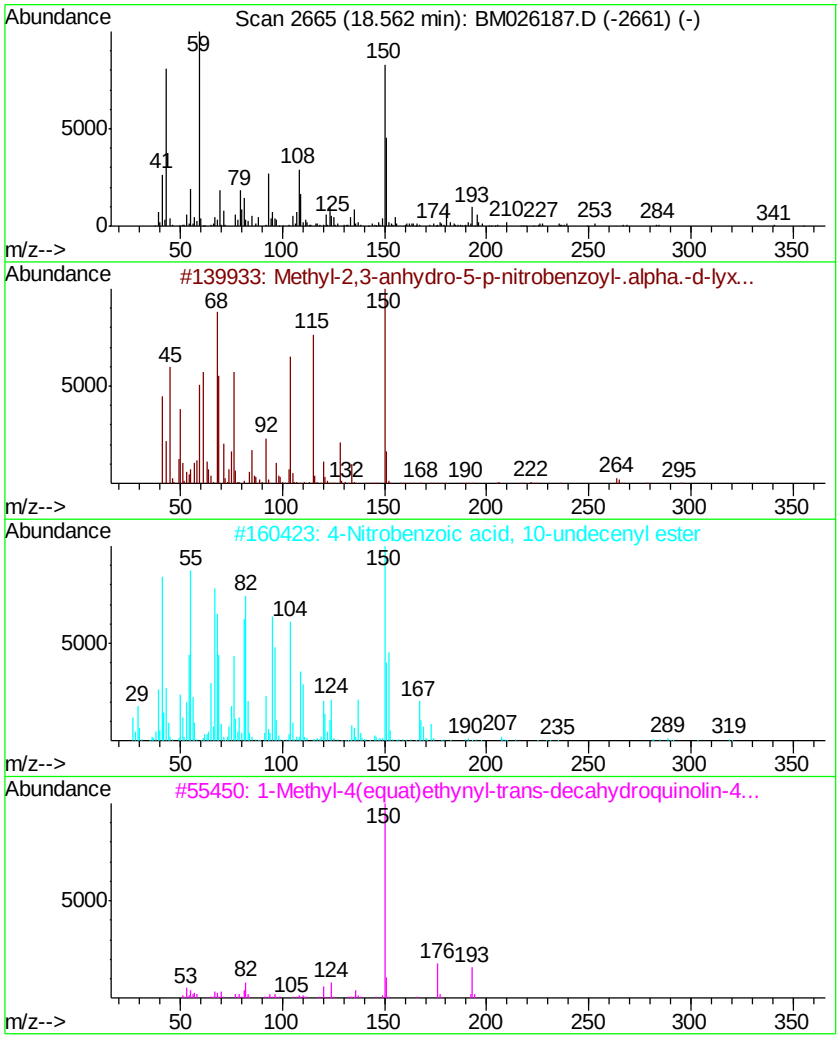
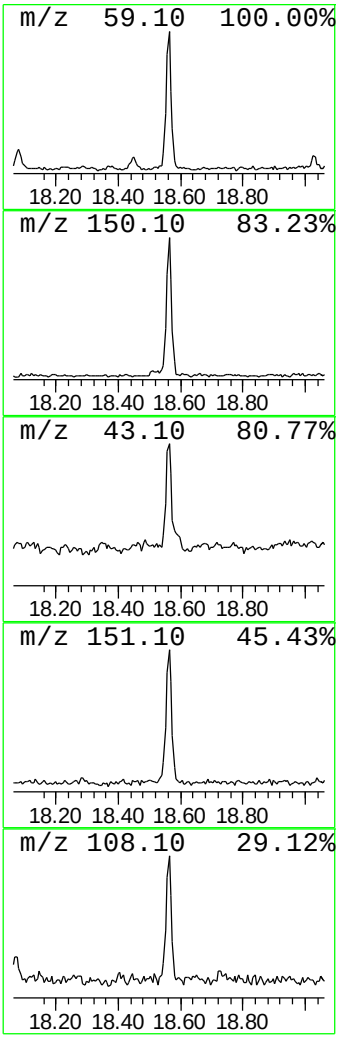
Peak Number 27 unknown-04 Concentration Rank 26

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.56	2.52 ng/ul	266574	Phenanthrene-d10	16.87

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Methyl-2,3-anhydro-5-p-nitrobenz...	295	C13H13NO7	1000214-61-0	43
2	4-Nitrobenzoic acid, 10-undecenyl...	319	C18H25NO4	1000281-86-3	27
3	1-Methyl-4(equat)ethynyl-trans-d...	193	C12H19NO	057581-44-9	22
4	4-Nitrobenzoic acid, tridec-2-yn...	345	C20H27NO4	1000299-27-3	22
5	7-Octen-2-ol, 2-methyl-6-methylene-	154	C10H18O	000543-39-5	22





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Acq On : 30 May 2020 00:46
Operator : MA/SJ
Sample : L2820-08
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : CB638

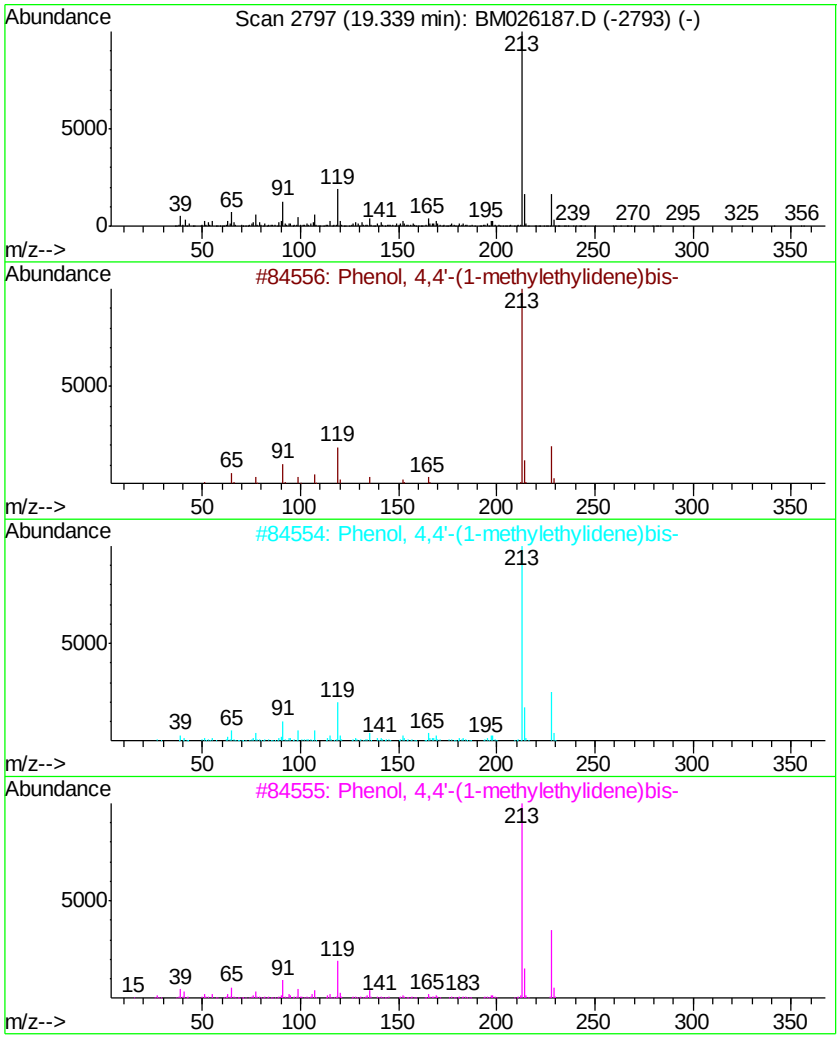
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM051320MA.M
Quant Title : SVOA CALIBRATION

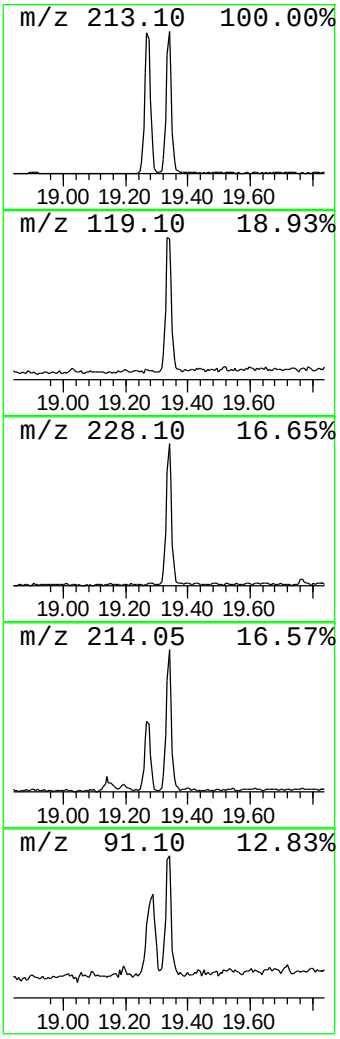
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 28 Phenol, 4,4'-(1-methylethyl... Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.34	4.64 ng/ul	563912	Chrysene-d12	21.07

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, 4,4'-(1-methylethylidene...	228	C15H16O2	000080-05-7	97
2	Phenol, 4,4'-(1-methylethylidene...	228	C15H16O2	000080-05-7	93
3	Phenol, 4,4'-(1-methylethylidene...	228	C15H16O2	000080-05-7	64
4	2H,8H-Benzo[1,2-b:3,4-b']dipyran...	228	C14H12O3	000523-59-1	59
5	Carbanilic acid, p-phenyl-	213	C13H11NO2	004474-53-7	50





Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM052920\
Data File : BM026187.D
Acq On : 30 May 2020 00:46
Operator : MA/SJ
Sample : L2820-08
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : CB638

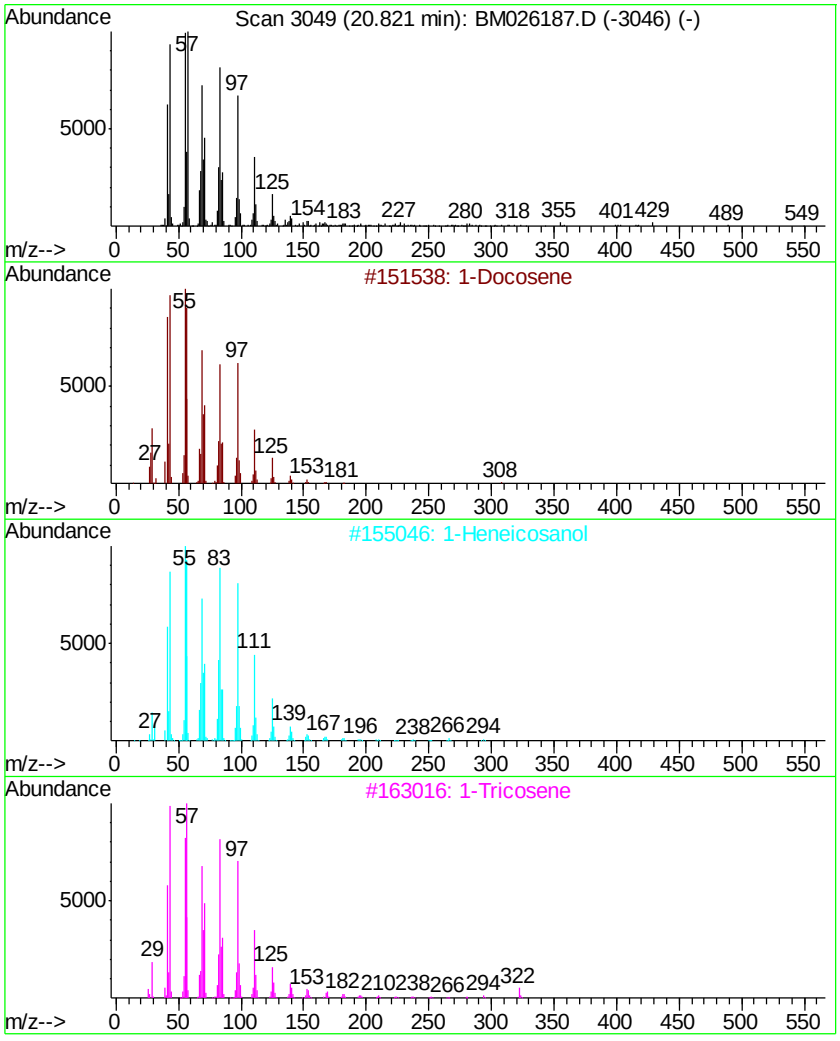
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM051320MA.M
Quant Title : SVOA CALIBRATION

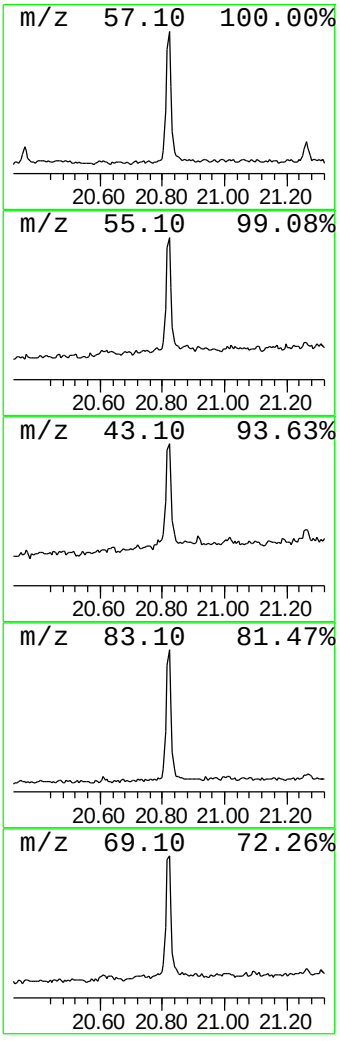
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 29 (DEL) Alkane: Straight-Chai... Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.82	3.99 ng/ul	483984	Chrysene-d12	21.07

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Docosene	308	C22H44	001599-67-3	94
2	1-Heneicosanol	312	C21H44O	015594-90-8	93
3	1-Tricosene	322	C23H46	018835-32-0	91
4	Octacosanol	410	C28H58O	000557-61-9	91
5	1-Heneicosyl formate	340	C22H44O2	077899-03-7	91





Tentatively Identified Compound (LSC) summary

Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM052920\
Data File : BM026187.D
Acq On : 30 May 2020 00:46
Operator : MA/SJ
Sample : L2820-08
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : CB638

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
Paraldehyde	3.84	2.6	ng/ul	122966	1	7.47	941387	20.0
1,3-Oxathiolane	4.18	6.0	ng/ul	280910	1	7.47	941387	20.0
Bicyclo[3.1.1]hep...	6.18	2.1	ng/ul	100999	1	7.47	941387	20.0
Bicyclo[2.2.1]hep...	6.48	5.0	ng/ul	234027	1	7.47	941387	20.0
p-Cymene	7.62	72.2	ng/ul	3398970	1	7.47	941387	20.0
Benzenamine, 3-me...	8.46	163.1	ng/ul	7676540	1	7.47	941387	20.0
Bicyclo[2.2.1]hep...	8.70	41.9	ng/ul	1974190	1	7.47	941387	20.0
Hexanoic acid, 3,...	9.10	3.6	ng/ul	216906	2	10.23	1190080	20.0
1,6-Octadiene, 5,...	9.51	3.1	ng/ul	187654	2	10.23	1190080	20.0
Cyclohexanemethan...	9.62	96.7	ng/ul	5755920	2	10.23	1190080	20.0
Bicyclo[2.2.1]hep...	9.66	54.8	ng/ul	3259610	2	10.23	1190080	20.0
unknown-01	9.79	6.5	ng/ul	385074	2	10.23	1190080	20.0
Terpinen-4-ol	10.12	10.6	ng/ul	629469	2	10.23	1190080	20.0
.alpha.-Terpineol	10.30	3.4	ng/ul	199206	2	10.23	1190080	20.0
Phenol, 2-propyl-	11.08	18.6	ng/ul	1109230	2	10.23	1190080	20.0
Phenol, p-tert-bu...	11.60	3.8	ng/ul	227256	2	10.23	1190080	20.0
Benzenamine, 3-ch...	11.75	46.3	ng/ul	2756760	2	10.23	1190080	20.0
Benzenamine, 3-ch...	11.86	9.1	ng/ul	543768	2	10.23	1190080	20.0
Cyclohexanemethan...	11.99	3.1	ng/ul	185857	2	10.23	1190080	20.0
unknown-02	14.05	2.3	ng/ul	193817	3	14.12	1672800	20.0
1-Naphthalenol, 2...	15.38	2.0	ng/ul	171830	3	14.12	1672800	20.0
Benzenesulfonamid...	15.65	14.0	ng/ul	1478910	4	16.87	2119510	20.0
2(3H)-Benzothiazo...	15.82	5.2	ng/ul	549868	4	16.87	2119510	20.0
Benzenesulfonamid...	16.16	8.9	ng/ul	941565	4	16.87	2119510	20.0
unknown-03	16.48	9.4	ng/ul	995673	4	16.87	2119510	20.0
Phenol, 3-(2-phen...	17.07	4.3	ng/ul	451034	4	16.87	2119510	20.0
unknown-04	18.56	2.5	ng/ul	266574	4	16.87	2119510	20.0
Phenol, 4,4'-(1-m...	19.34	4.6	ng/ul	563912	5	21.07	2428890	20.0
(DEL) Alkane: Str...	20.82	4.0	ng/ul	483984	5	21.07	2428890	20.0