LSC Area Percent Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM062320\
Data File : BM026527.D
Acq On : 24 Jun 2020 09:33
Operator : JU/CG
Sample : L3069-11DL 2X
Misc :
ALS Vial : 37 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : CB7H0DL

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: BM026530.D

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.993	15	18	32	rVB	185944	270738	12.62%	1.179%
2	3.663	128	132	137	rBV2	7164	11405	0.53%	0.050%
3	3.746	142	146	151	rVV	56243	74379	3.47%	0.324%
4	4.087	199	204	217	rVB	53355	92185	4.30%	0.402%
5	5.069	367	371	379	rVB4	7420	14814	0.69%	0.065%
6	5.351	415	419	426	rVB9	5947	12845	0.60%	0.056%
7	5.428	426	432	437	rBV6	6722	12905	0.60%	0.056%
8	6.369	588	592	600	rVB4	12011	24619	1.15%	0.107%
9	6.592	622	630	643	rVB3	41012	103627	4.83%	0.451%
10	6.734	649	654	665	rBV	131529	213794	9.97%	0.931%
11	6.869	672	677	680	rBV4	8065	12381	0.58%	0.054%
12	6.922	680	686	694	rVV	156136	256031	11.94%	1.115%
13	7.375	756	763	772	rVV	440857	691441	32.24%	3.012%
14	7.463	772	778	784	rVV3	19796	47622	2.22%	0.207%
15	7.516	784	787	795	rVB	23156	34256	1.60%	0.149%
16	7.745	822	826	830	rVB5	8283	13473	0.63%	0.059%
17	8.104	881	887	894	rVV	125301	215576	10.05%	0.939%
18	8.157	894	896	904	rVB7	11202	21562	1.01%	0.094%
19	8.357	923	930	940	rBV	1385829	2081757	97.06%	9.068%
20	8.475	946	950	953	rBV	20537	32332	1.51%	0.141%
21	8.522	953	958	966	rVV	176918	297821	13.89%	1.297%
22	8.604	966	972	980	rVB	51847	82846	3.86%	0.361%
23	8.716	986	991	999	rVB6	12778	27528	1.28%	0.120%
24	8.875	1011	1018	1022	rBV4	20164	35348	1.65%	0.154%
25	9.010	1032	1041	1047	rBV2	41692	82295	3.84%	0.358%
26	9.239	1073	1080	1093	rBV	151980	274056	12.78%	1.194%
27	9.363	1096	1101	1106	rVV5	17848	31657	1.48%	0.138%
28	9.416	1106	1110	1115	rVV3	25898	43432	2.02%	0.189%
29	9.457	1115	1117	1122	rVB5	7999	10602	0.49%	0.046%
30	9.522	1122	1128	1131	rBV	91736	143349	6.68%	0.624%
31	9.557	1131	1134	1142	rVB2	87437	136787	6.38%	0.596%
32	9.686	1148	1156	1166	rBV2	47265	102898	4.80%	0.448%
33	9.780	1166	1172	1177	rBV	219448	396569	18.49%	1.727%
34	9.945	1197	1200	1206	rVV6	12025	24361	1.14%	0.106%

LSC Area Percent Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM062320\
Data File : BM026527.D
Acq On : 24 Jun 2020 09:33
Operator : JU/CG
Sample : L3069-11DL 2X
Misc :
ALS Vial : 37 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : CB7H0DL

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	10.133	1224	1232	1238	rVV	661344	1079560	50.33%	4.703%
36	10.180	1238	1240	1250	rVB4	19662	43637	2.03%	0.190%
37	10.286	1250	1258	1270	rBV	205367	435615	20.31%	1.898%
38	10.386	1272	1275	1279	rVV6	5959	10608	0.49%	0.046%
39	10.439	1279	1284	1291	rVV6	12923	28366	1.32%	0.124%
40	10.598	1307	1311	1322	rBV9	9283	24535	1.14%	0.107%
41	10.727	1329	1333	1339	rVV7	10482	23473	1.09%	0.102%
42	10.792	1340	1344	1348	rVV7	11615	22525	1.05%	0.098%
43	10.845	1348	1353	1358	rVB7	12094	26616	1.24%	0.116%
44	10.992	1371	1378	1389	rVV6	24456	86425	4.03%	0.376%
45	11.139	1398	1403	1407	rVB8	9399	18752	0.87%	0.082%
46	11.210	1411	1415	1420	rVB6	10752	17703	0.83%	0.077%
47	11.345	1434	1438	1448	rVB4	20264	41129	1.92%	0.179%
48	11.510	1459	1466	1471	rVB2	35955	63100	2.94%	0.275%
49	11.657	1484	1491	1499	rBV	440009	727248	33.91%	3.168%
50	11.727	1499	1503	1505	rVV5	16456	31240	1.46%	0.136%
51	11.769	1505	1510	1522	rVV	84963	180739	8.43%	0.787%
52	11.886	1524	1530	1537	rVB9	13267	33274	1.55%	0.145%
53	12.174	1575	1579	1587	rVB3	16565	27483	1.28%	0.120%
54	12.380	1610	1614	1620	rBV8	8961	17679	0.82%	0.077%
55	12.680	1662	1665	1669	rBV5	5950	10419	0.49%	0.045%
56	12.751	1672	1677	1680	rVV6	7788	15345	0.72%	0.067%
57	12.833	1686	1691	1696	rBV8	15191	25762	1.20%	0.112%
58	12.886	1696	1700	1705	rVB2	22186	33076	1.54%	0.144%
59	12.951	1706	1711	1721	rBV2	35881	59605	2.78%	0.260%
60	13.080	1729	1733	1740	rVB2	25182	38680	1.80%	0.168%
61	13.410	1786	1789	1792	rVV4	8189	11971	0.56%	0.052%
62	13.463	1792	1798	1802	rVV	414154	569371	26.55%	2.480%
63	13.510	1802	1806	1813	rVB	161825	219761	10.25%	0.957%
64	13.592	1818	1820	1825	rVV6	9632	14579	0.68%	0.064%
65	13.639	1825	1828	1836	rVB7	14154	22820	1.06%	0.099%
66	13.721	1836	1842	1851	rBV	395715	598667	27.91%	2.608%
67	13.963	1879	1883	1888	rVB2	53076	76008	3.54%	0.331%
68	14.033	1889	1895	1903	rBV	1059083	1483117	69.15%	6.461%
69	14.098	1904	1906	1909	rVB3	11316	11096	0.52%	0.048%
70	14.321	1939	1944	1948	rBV7	17979	35043	1.63%	0.153%
71	14.657	1994	2001	2009	rBV7	11348	28575	1.33%	0.124%

LSC Area Percent Report

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Acq On : 24 Jun 2020 09:33
Operator : JU/CG
Sample : L3069-11DL 2X
Misc :
ALS Vial : 37 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : CB7H0DL

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	14.751	2012	2017	2023	rBV2	18145	36937	1.72%	0.161%
73	14.810	2023	2027	2031	rBV	20522	32070	1.50%	0.140%
74	15.039	2060	2066	2072	rBV	512180	699693	32.62%	3.048%
75	15.092	2072	2075	2078	rVB5	9173	11117	0.52%	0.048%
76	15.186	2086	2091	2104	rBV2	112245	199314	9.29%	0.868%
77	15.309	2108	2112	2116	rVV2	28384	35984	1.68%	0.157%
78	15.509	2143	2146	2151	rVB5	12803	15064	0.70%	0.066%
79	15.568	2151	2156	2165	rBV	323033	439907	20.51%	1.916%
80	15.751	2181	2187	2192	rBV3	60408	120948	5.64%	0.527%
81	15.909	2210	2214	2218	rBV2	16352	19585	0.91%	0.085%
82	15.951	2218	2221	2223	rBV2	10936	10695	0.50%	0.047%
83	16.086	2239	2244	2251	rBV	205377	274732	12.81%	1.197%
84	16.151	2252	2255	2259	rVB4	15797	20301	0.95%	0.088%
85	16.368	2288	2292	2294	rBV2	31430	45489	2.12%	0.198%
86	16.586	2327	2329	2332	rBV4	8110	11788	0.55%	0.051%
87	16.668	2341	2343	2350	rVB7	10278	12419	0.58%	0.054%
88	16.792	2358	2364	2369	rBV	1319386	1795990	83.74%	7.823%
89	16.886	2375	2380	2391	rVB	545314	787432	36.71%	3.430%
90	17.215	2432	2436	2443	rVB7	22955	39301	1.83%	0.171%
91	17.727	2519	2523	2526	rBV3	28515	44213	2.06%	0.193%
92	17.998	2565	2569	2579	rBV7	38101	69200	3.23%	0.301%
93	18.486	2648	2652	2656	rVB2	65746	79974	3.73%	0.348%
94	18.662	2680	2682	2688	rBV6	11756	18177	0.85%	0.079%
95	19.197	2768	2773	2780	rBV	735455	940963	43.87%	4.099%
96	19.274	2782	2786	2791	rVB	83537	112723	5.26%	0.491%
97	20.756	3034	3038	3045	rBV	316775	360816	16.82%	1.572%
98	21.003	3075	3080	3090	rBV	1676118	2144847	100.00%	9.343%
99	22.962	3408	3413	3420	rBV	344117	544559	25.39%	2.372%
100	23.091	3430	3435	3446	rVB	1098677	1911484	89.12%	8.327%

Sum of corrected areas: 22956615

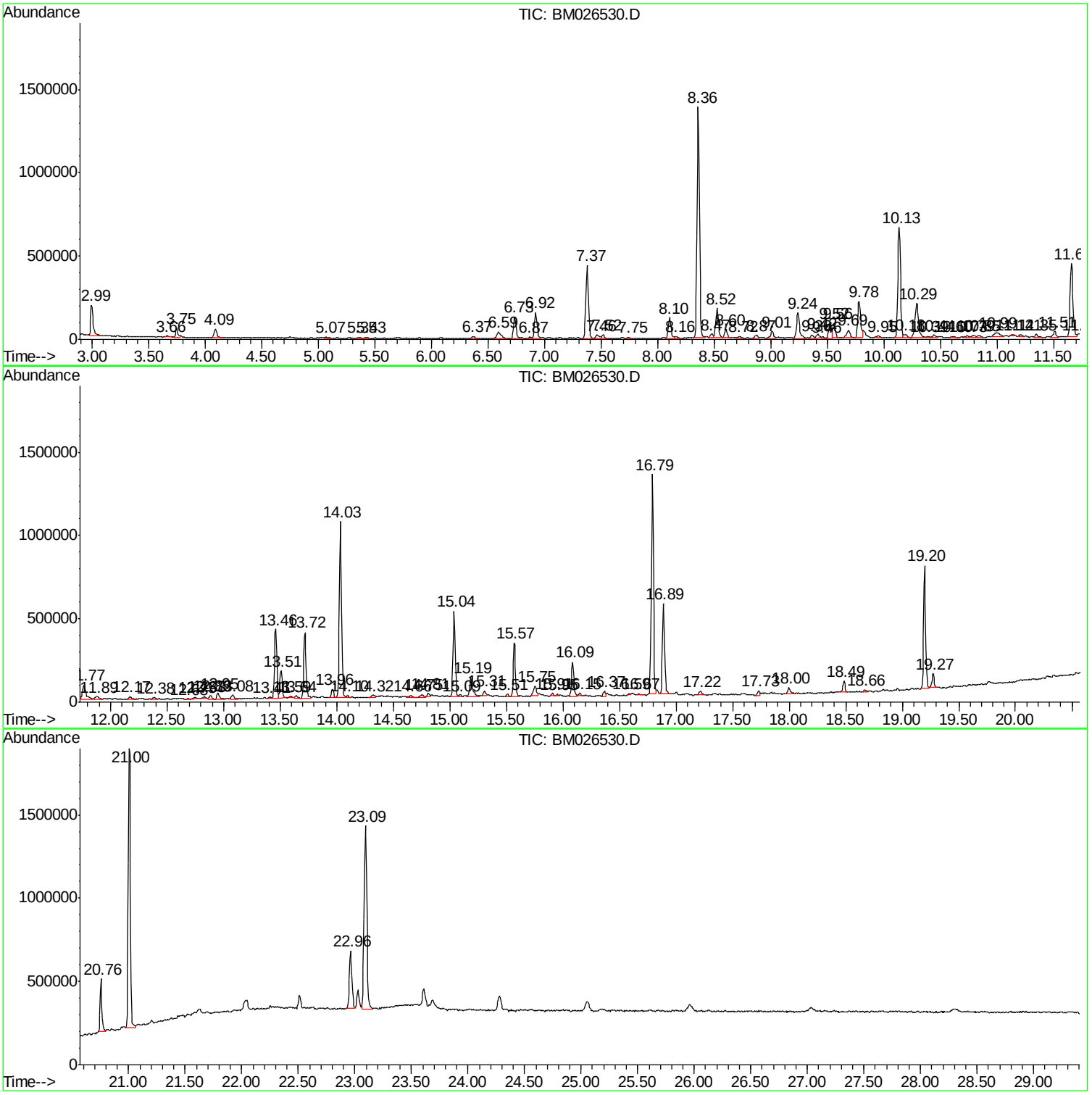
LSC Report - Integrated Chromatogram

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM062320\
Data File : BM026527.D
Acq On : 24 Jun 2020 09:33
Operator : JU/CG
Sample : L3069-11DL 2X
Misc :
ALS Vial : 37 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : CB7H0DL

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 : BM026527.D
Acq On : 24 Jun 2020 09:33
Operator : JU/CG
Sample : L3069-11DL 2X
Misc :
ALS Vial : 37 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : CB7H0DL

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

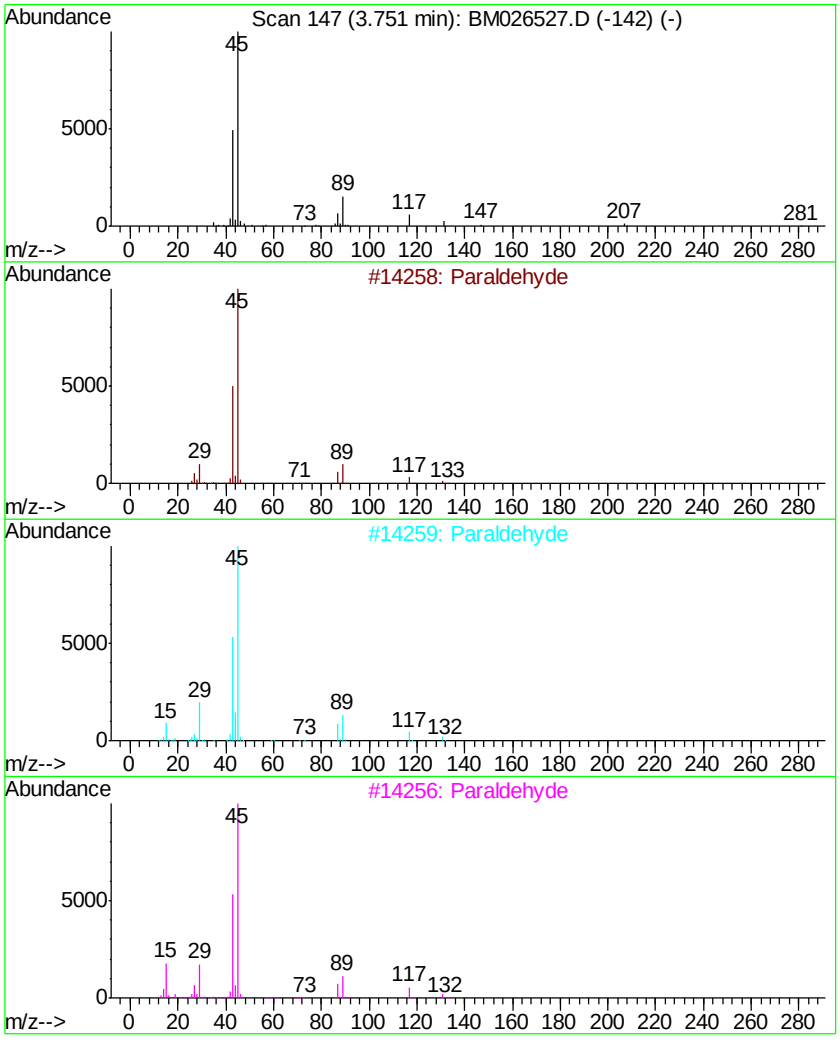
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

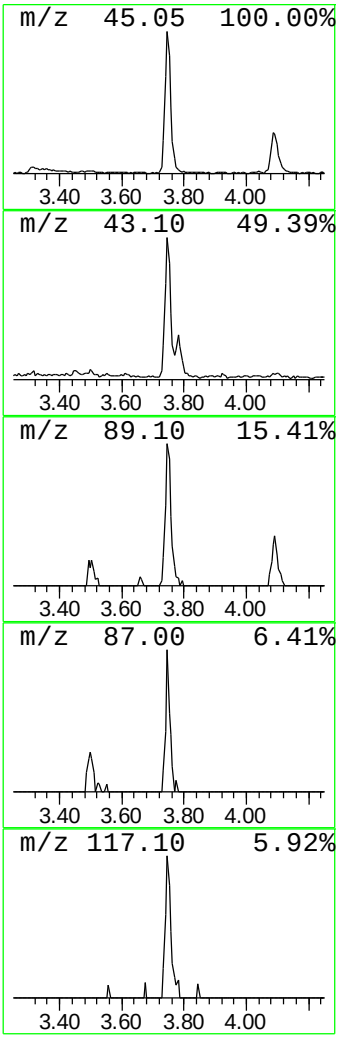
****

Peak Number 1 Paraldehyde Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.75	2.15 ng/ul	74379	1,4-Dichlorobenzene-d4	7.37

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	56
4	Ethyl 3-methylbutan-2-yl carbonate	160	C8H16O3	1000373-78-2	50
5	Propane, 2,2'-[ethylidenebis(oxy...	146	C8H18O2	004285-59-0	50





Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM062320\
Data File : BM026527.D
Acq On : 24 Jun 2020 09:33
Operator : JU/CG
Sample : L3069-11DL 2X
Misc :
ALS Vial : 37 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : CB7H0DL

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

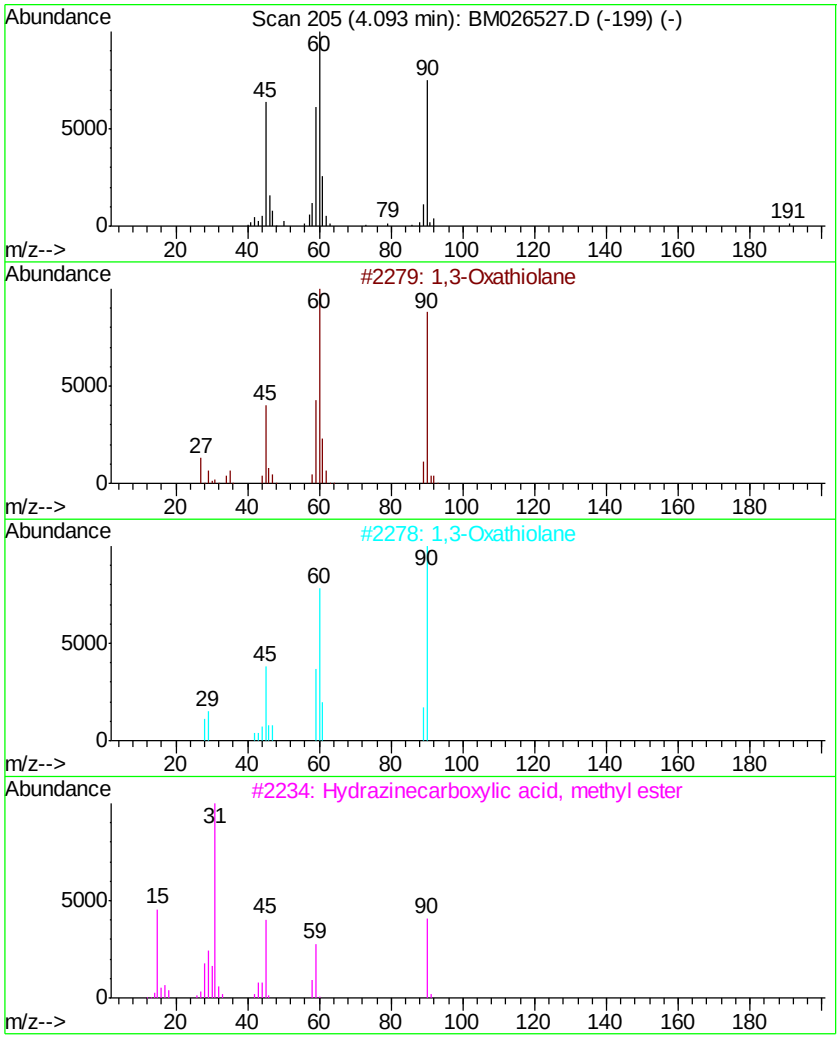
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

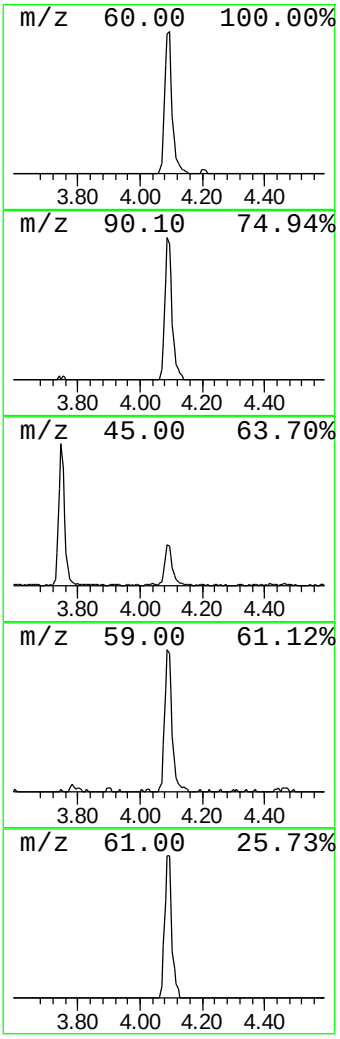
****

Peak Number 2 1,3-Oxathiolane Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.09	2.67 ng/ul	92185	1,4-Dichlorobenzene-d4	7.37

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	Hydrazinecarboxylic acid, methyl...	90	C2H6N2O2	006294-89-9	59
4	Acrolein, 2-chloro-	90	C3H3ClO	000683-51-2	59
5	1-Cyano-N-fluoroformimidoyl fluo...	90	C2F2N2	014011-05-3	40





Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM062320\
Data File : BM026527.D
Acq On : 24 Jun 2020 09:33
Operator : JU/CG
Sample : L3069-11DL 2X
Misc :
ALS Vial : 37 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : CB7H0DL

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

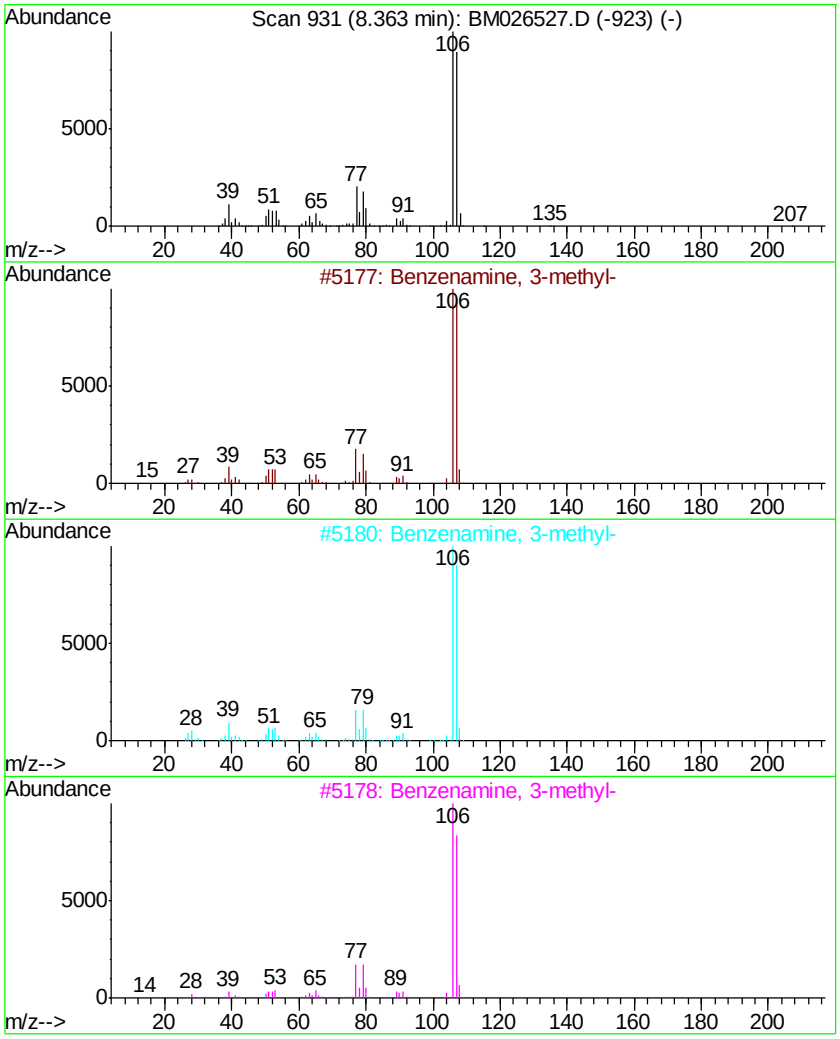
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

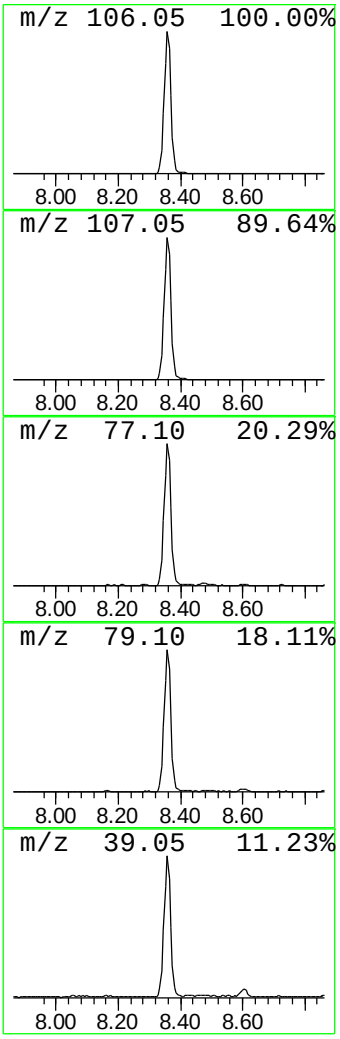
****

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.36	60.22 ng/ul	2081760	1,4-Dichlorobenzene-d4	7.37

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzenamine, 3-methyl-	107	C7H9N	000108-44-1	97
2	Benzenamine, 3-methyl-	107	C7H9N	000108-44-1	97
3	Benzenamine, 3-methyl-	107	C7H9N	000108-44-1	97
4	Benzenamine, 3-methyl-	107	C7H9N	000108-44-1	96
5	o-Toluidine	107	C7H9N	000095-53-4	93





Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM062320\
Data File : BM026527.D
Acq On : 24 Jun 2020 09:33
Operator : JU/CG
Sample : L3069-11DL 2X
Misc :
ALS Vial : 37 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : CB7H0DL

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

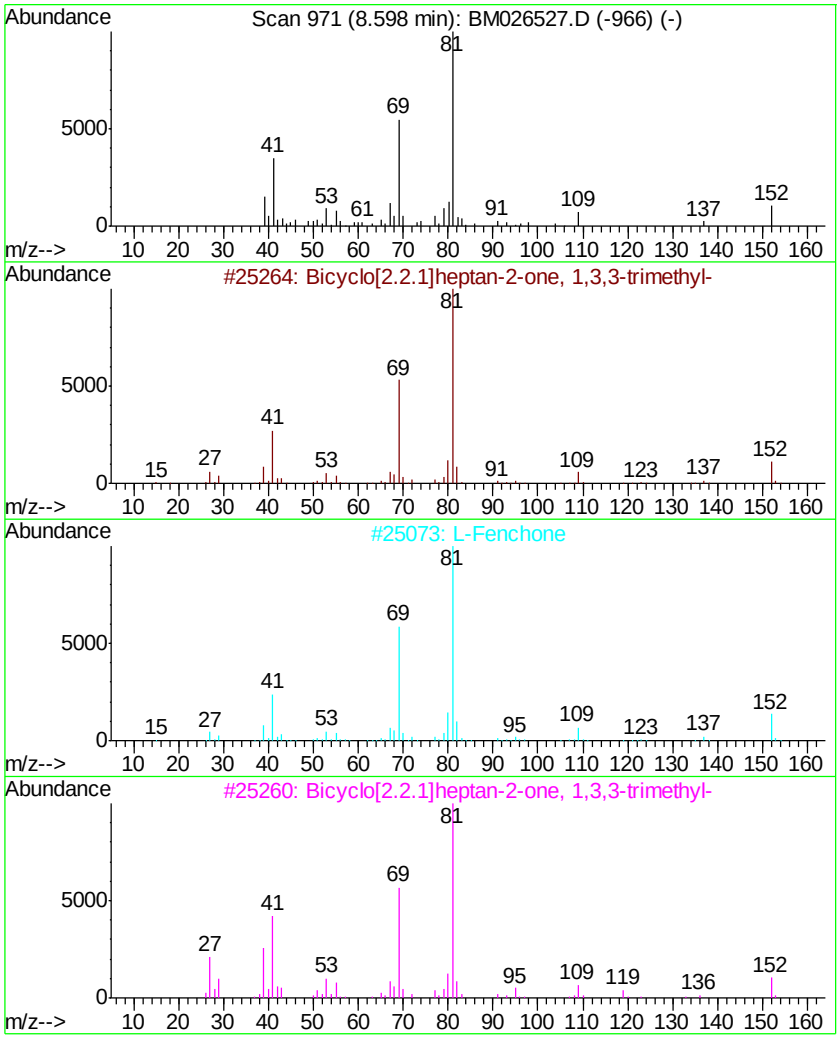
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

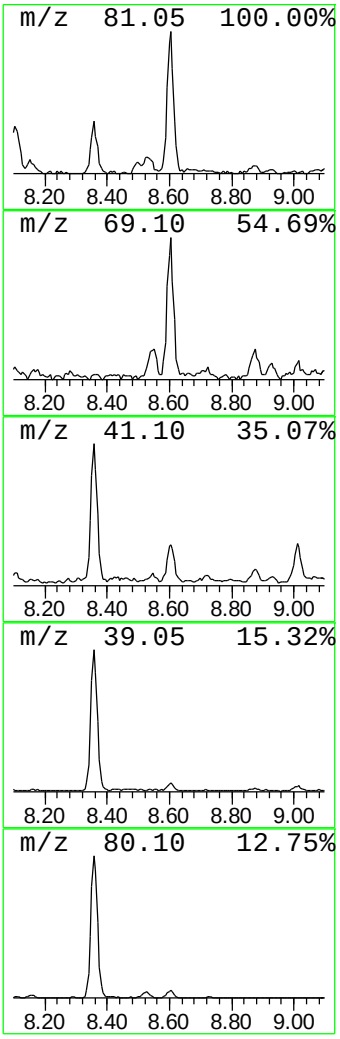
****

Peak Number 4 Bicyclo[2.2.1]heptan-2-one,... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.60	2.40 ng/ul	82846	1,4-Dichlorobenzene-d4	7.37

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Bicyclo[2.2.1]heptan-2-one, 1,3,...	152	C10H16O	001195-79-5	91
2	L-Fenchone	152	C10H16O	007787-20-4	87
3	Bicyclo[2.2.1]heptan-2-one, 1,3,...	152	C10H16O	001195-79-5	87
4	L-Fenchone	152	C10H16O	007787-20-4	87
5	Bicyclo[2.2.1]heptan-2-one, 1,3,...	152	C10H16O	001195-79-5	87





Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM062320\
Data File : BM026527.D
Acq On : 24 Jun 2020 09:33
Operator : JU/CG
Sample : L3069-11DL 2X
Misc :
ALS Vial : 37 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : CB7H0DL

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

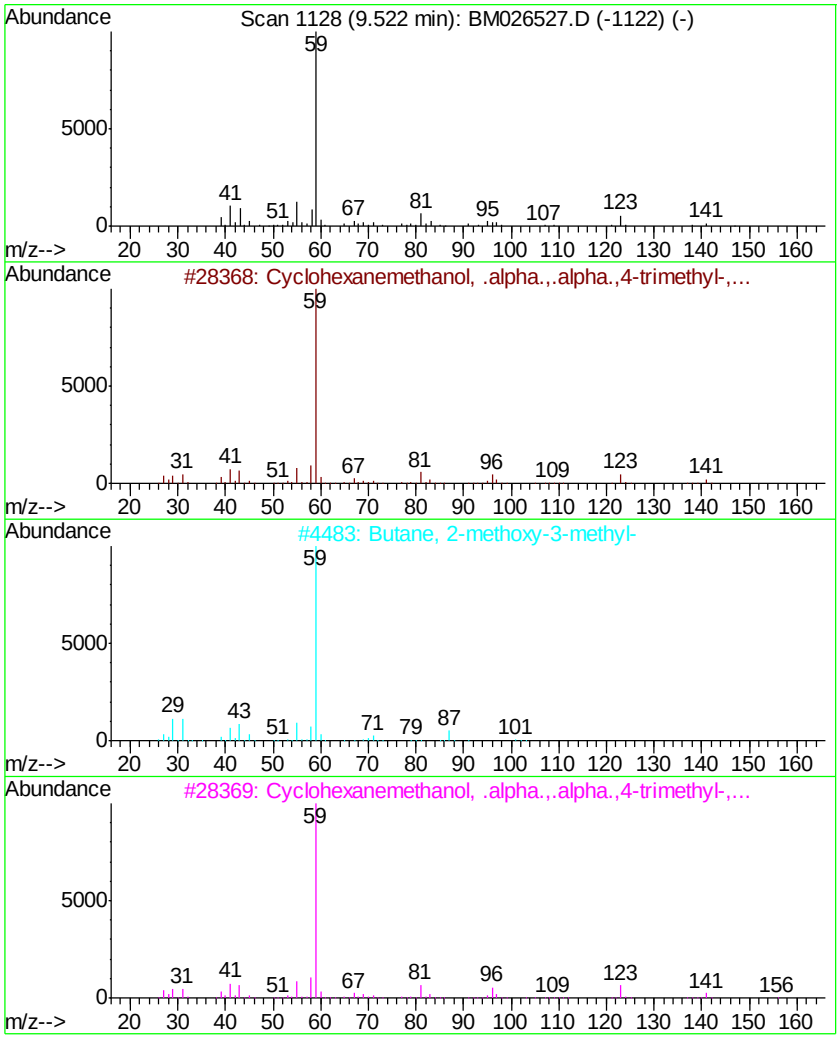
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

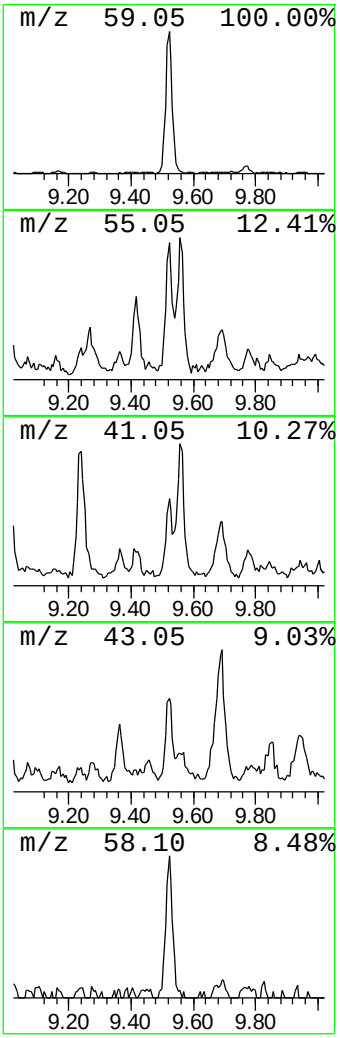
****

Peak Number 5 Cyclohexanemethanol, .alpha... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.52	2.66 ng/ul	143349	Naphthalene-d8	10.13

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclohexanemethanol, .alpha.,.al...	156	C10H20O	007322-63-6	78
2	Butane, 2-methoxy-3-methyl-	102	C6H14O	062016-49-3	64
3	Cyclohexanemethanol, .alpha.,.al...	156	C10H20O	005114-00-1	64
4	o-Menthan-8-ol	156	C10H20O	343855-44-7	56
5	2,3-Dimethyl-4-penten-2-ol	114	C7H14O	019781-52-3	50





Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM062320\
Data File : BM026527.D
Acq On : 24 Jun 2020 09:33
Operator : JU/CG
Sample : L3069-11DL 2X
Misc :
ALS Vial : 37 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : CB7H0DL

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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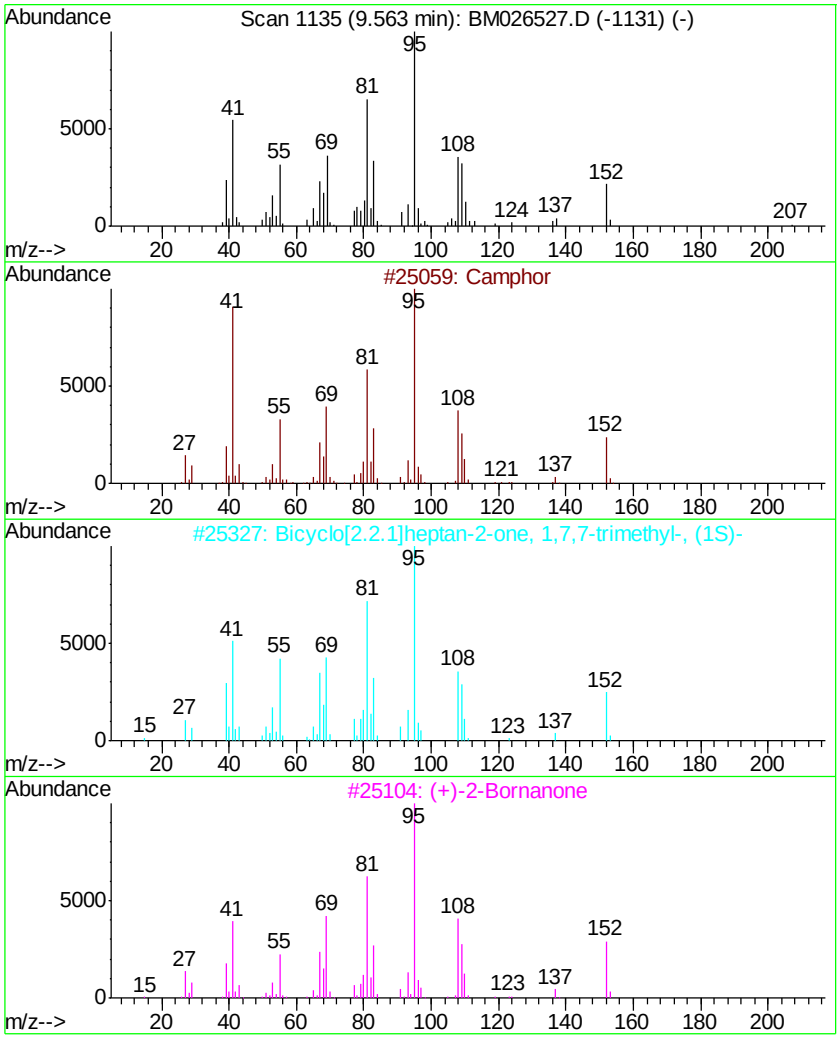
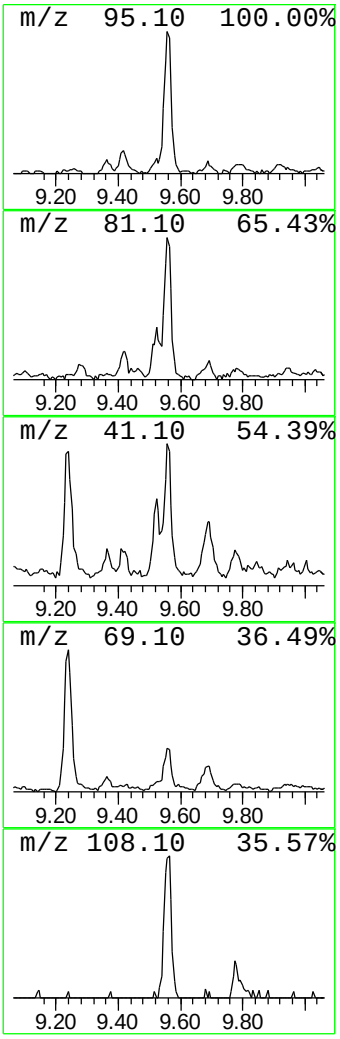
Peak Number 6 Camphor Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.56	2.53 ng/ul	136787	Naphthalene-d8	10.13

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





Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM062320\
Data File : BM026527.D
Acq On : 24 Jun 2020 09:33
Operator : JU/CG
Sample : L3069-11DL 2X
Misc :
ALS Vial : 37 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : CB7H0DL

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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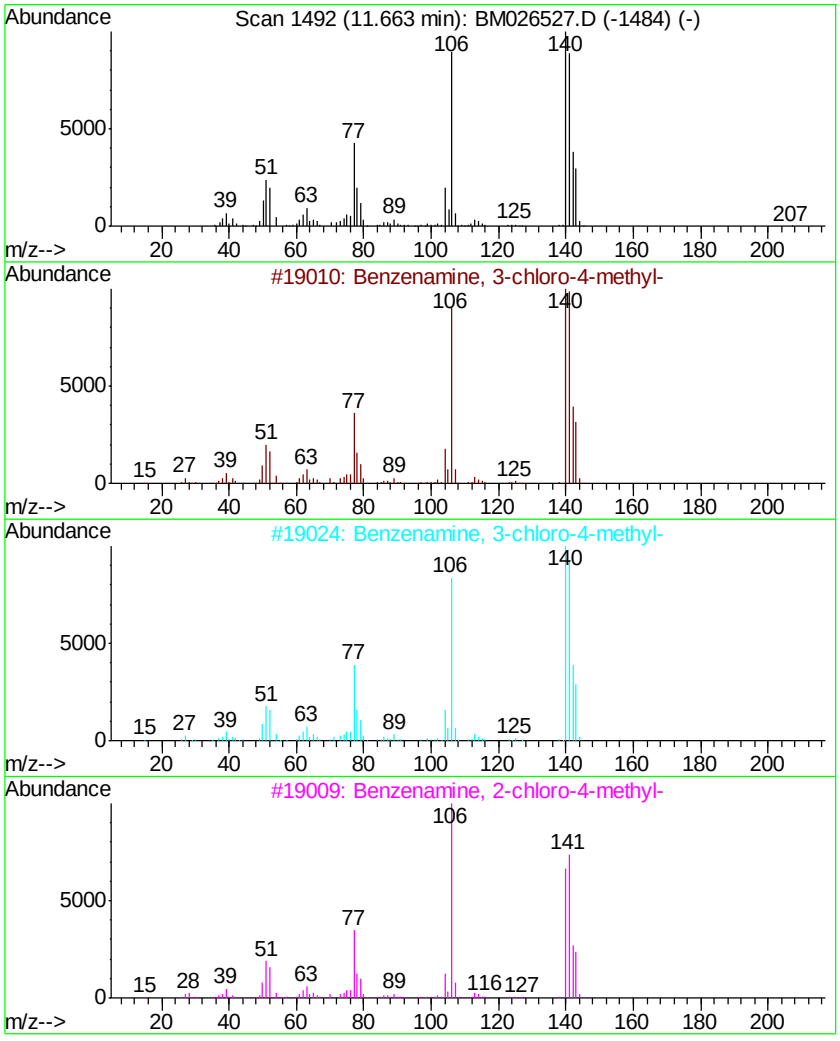
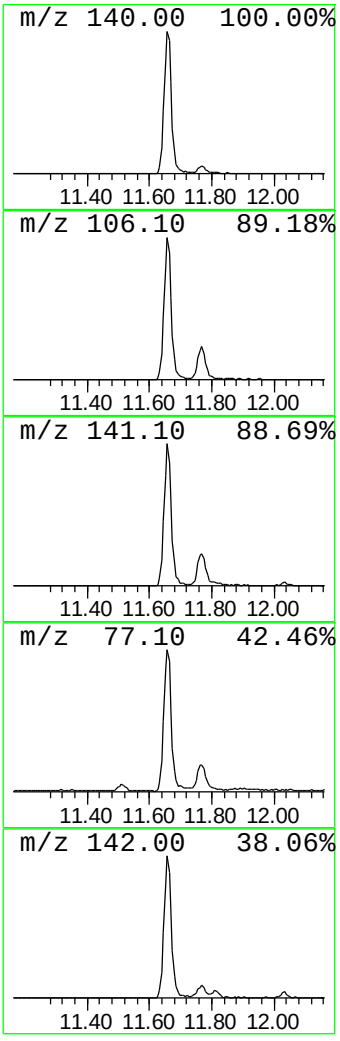
Peak Number 7 Benzenamine, 3-chloro-4-met... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.66	13.47 ng/ul	727248	Naphthalene-d8	10.13

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	94
5	Benzenamine, 3-chloro-4-methyl-	141	C7H8ClN	000095-74-9	90





Library Search Compound Report

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Data File : BM026527.D
Acq On : 24 Jun 2020 09:33
Operator : JU/CG
Sample : L3069-11DL 2X
Misc :
ALS Vial : 37 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : CB7H0DL

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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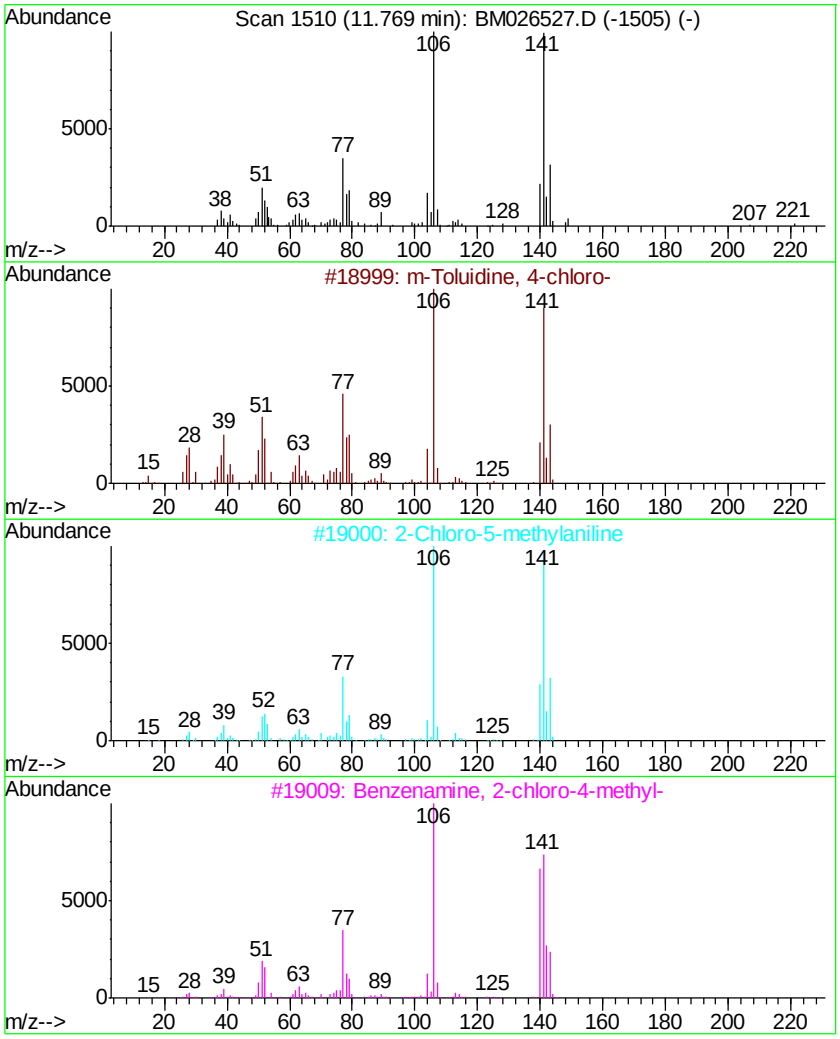
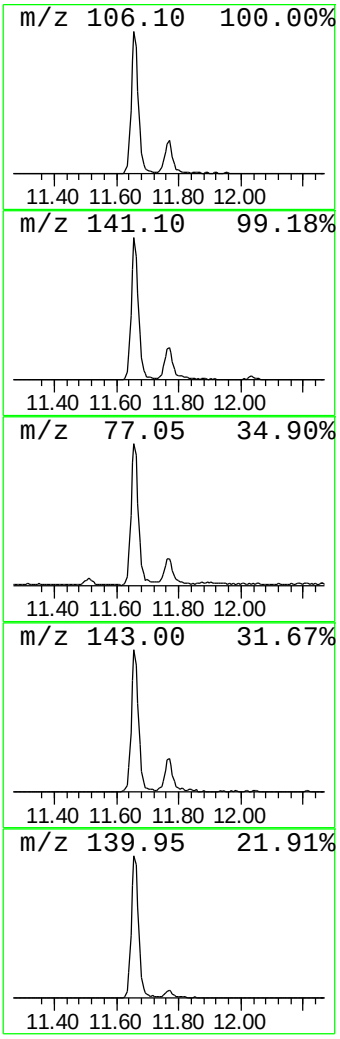
Peak Number 8 m-Toluidine, 4-chloro- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.77	3.35 ng/ul	180739	Naphthalene-d8	10.13

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	m-Toluidine, 4-chloro-	141	C7H8ClN	007149-75-9	96
2	2-Chloro-5-methylaniline	141	C7H8ClN	000095-81-8	90
3	Benzenamine, 2-chloro-4-methyl-	141	C7H8ClN	000615-65-6	87
4	Benzenamine, 2-chloro-6-methyl-	141	C7H8ClN	000087-63-8	87
5	Benzenamine, 5-chloro-2-methyl-	141	C7H8ClN	000095-79-4	87





Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM062320\
Data File : BM026527.D
Acq On : 24 Jun 2020 09:33
Operator : JU/CG
Sample : L3069-11DL 2X
Misc :
ALS Vial : 37 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : CB7H0DL

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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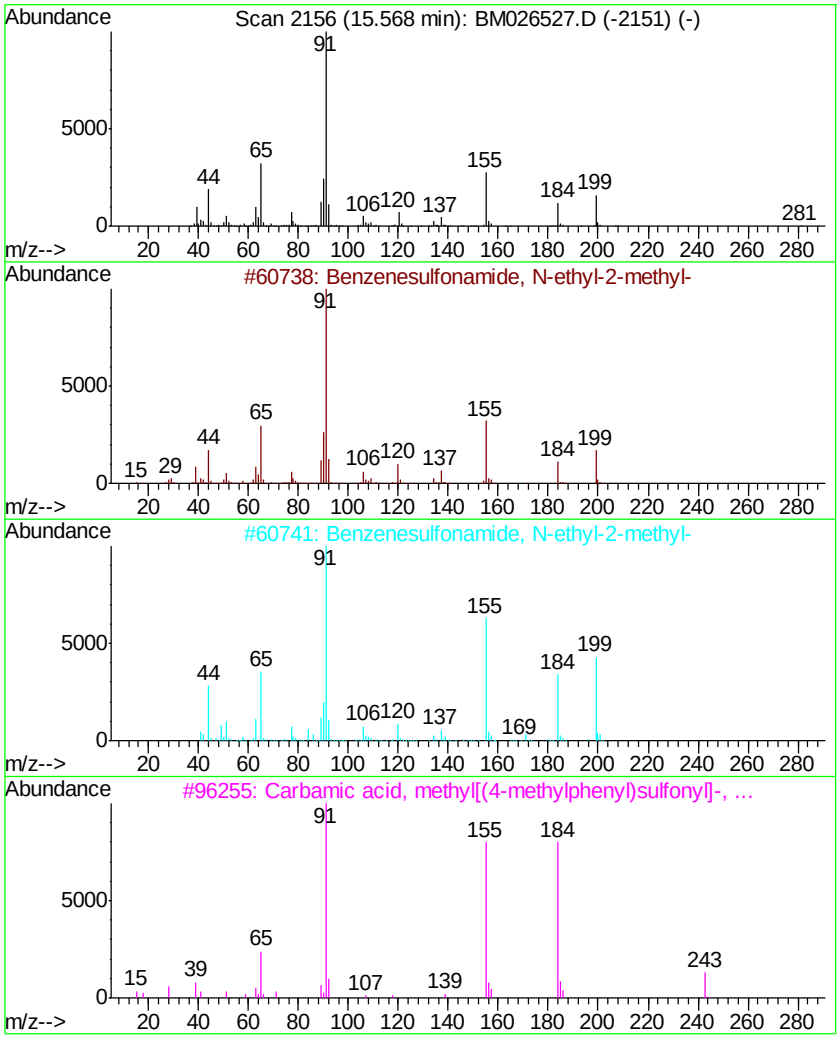
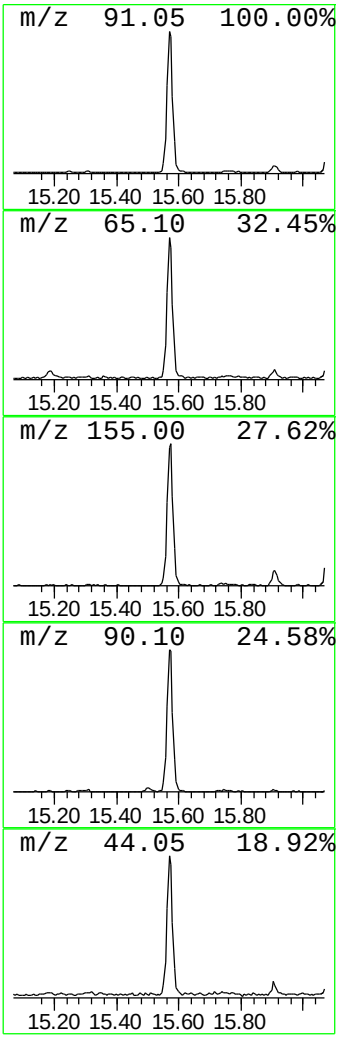
Peak Number 9 Benzenesulfonamide, N-ethyl... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.57	4.90 ng/ul	439907	Phenanthrene-d10	16.79

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzenesulfonamide, N-ethyl-2-me...	199	C9H13NO2S	001077-56-1	97
2	Benzenesulfonamide, N-ethyl-2-me...	199	C9H13NO2S	001077-56-1	95
3	Carbamic acid, methyl[(4-methylp...	243	C10H13NO4S	032258-50-7	53
4	Benzenesulfonamide, N,N'-(dithio...	460	C18H24N2O4S4	023516-74-7	53
5	4-Toluenesulfonylmethyl isocyanide	195	C9H9NO2S	036635-61-7	50





Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM062320\
Data File : BM026527.D
Acq On : 24 Jun 2020 09:33
Operator : JU/CG
Sample : L3069-11DL 2X
Misc :
ALS Vial : 37 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : CB7H0DL

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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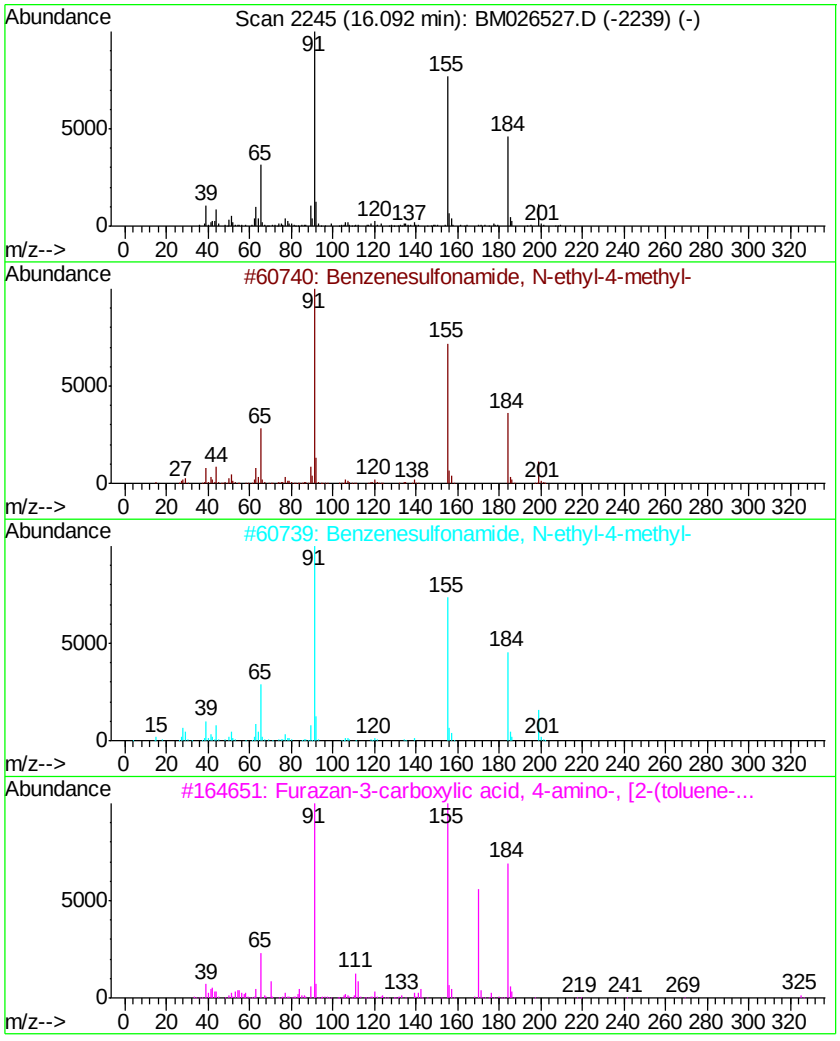
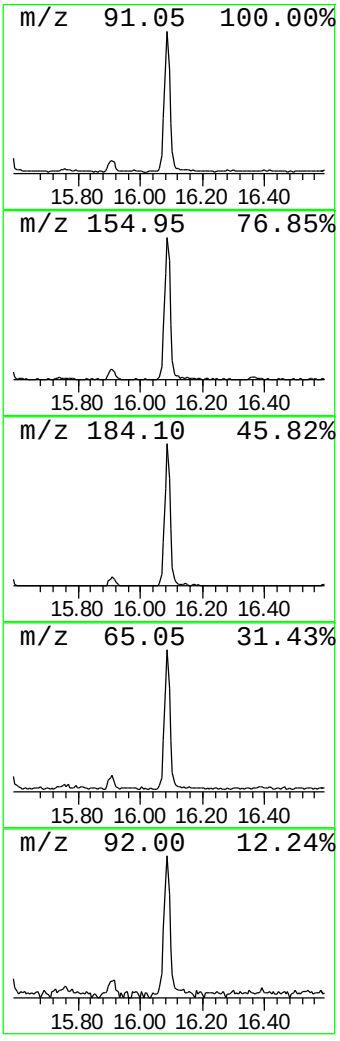
Peak Number 10 Benzenesulfonamide, N-ethyl... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.09	3.06 ng/ul	274732	Phenanthrene-d10	16.79

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzenesulfonamide, N-ethyl-4-me...	199	C9H13NO2S	000080-39-7	98
2	Benzenesulfonamide, N-ethyl-4-me...	199	C9H13NO2S	000080-39-7	93
3	Furazan-3-carboxylic acid, 4-ami...	325	C12H15N5O4S	1000304-69-4	78
4	4-Toluenesulfonylmethyl isocyanide	195	C9H9NO2S	036635-61-7	50
5	4-Toluenesulfonylmethyl isocyanide	195	C9H9NO2S	036635-61-7	50





Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM062320\
Data File : BM026527.D
Acq On : 24 Jun 2020 09:33
Operator : JU/CG
Sample : L3069-11DL 2X
Misc :
ALS Vial : 37 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : CB7H0DL

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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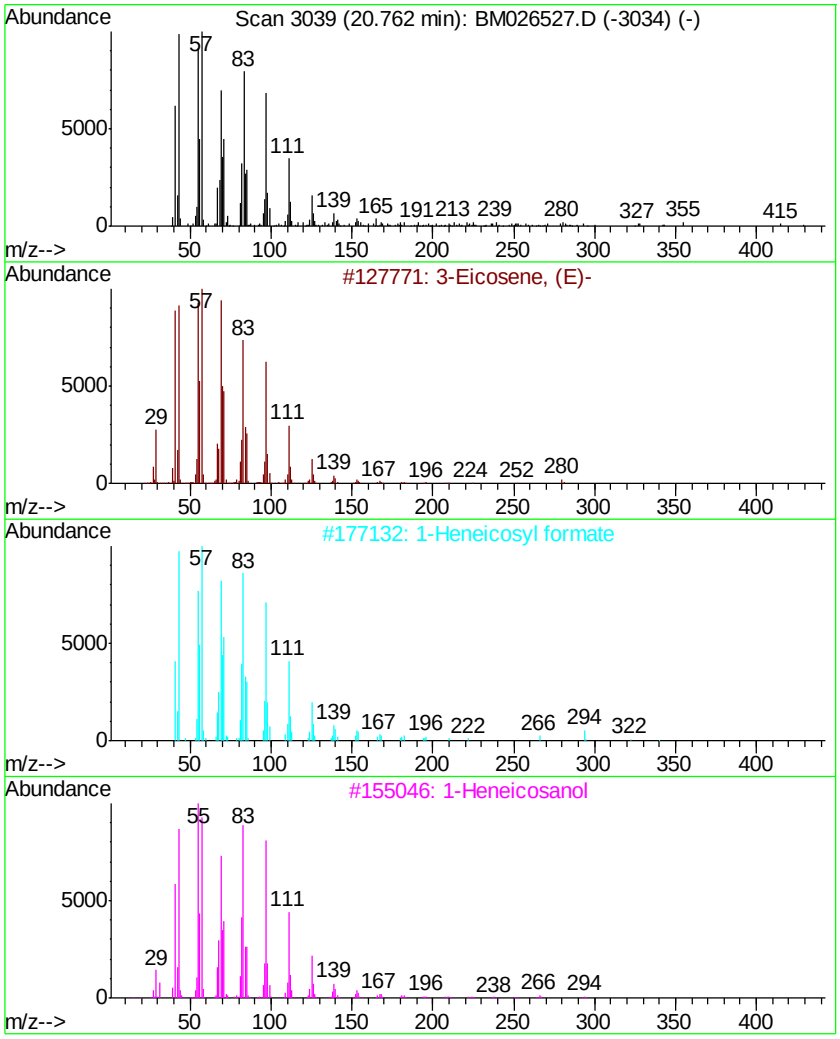
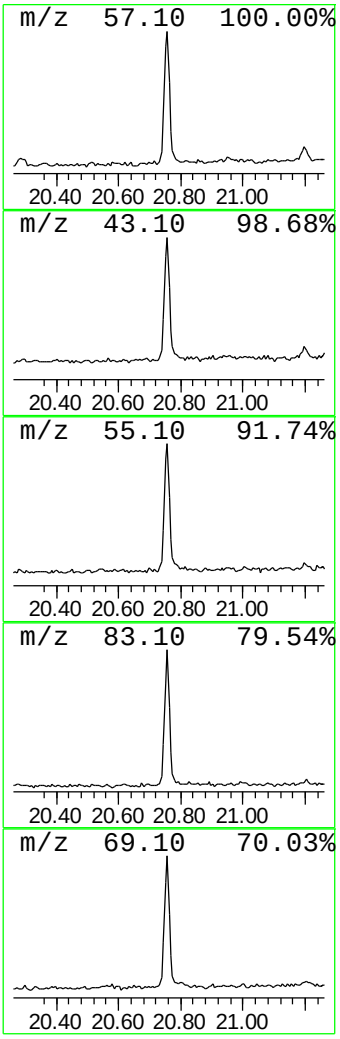
Peak Number 11 (DEL) Alkane: Straight-Chai... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.76	3.36 ng/ul	360816	Chrysene-d12	21.01

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Eicosene, (E)-	280	C20H40	074685-33-9	95
2	1-Heneicosyl formate	340	C22H44O2	077899-03-7	94
3	1-Heneicosanol	312	C21H44O	015594-90-8	94
4	n-Nonadecanol-1	284	C19H40O	001454-84-8	94
5	n-Nonadecanol-1	284	C19H40O	001454-84-8	91





Tentatively Identified Compound (LSC) summary

Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM062320\
Data File : BM026527.D
Acq On : 24 Jun 2020 09:33
Operator : JU/CG
Sample : L3069-11DL 2X
Misc :
ALS Vial : 37 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : CB7H0DL

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
Paraldehyde	3.75	2.1	ng/ul	74379	1	7.37	691441	20.0
1,3-Oxathiolane	4.09	2.7	ng/ul	92185	1	7.37	691441	20.0
Benzenamine, 3-me...	8.36	60.2	ng/ul	2081760	1	7.37	691441	20.0
Bicyclo[2.2.1]hep...	8.60	2.4	ng/ul	82846	1	7.37	691441	20.0
Cyclohexanemethan...	9.52	2.7	ng/ul	143349	2	10.13	1079560	20.0
Camphor	9.56	2.5	ng/ul	136787	2	10.13	1079560	20.0
Benzenamine, 3-ch...	11.66	13.5	ng/ul	727248	2	10.13	1079560	20.0
m-Toluidine, 4-ch...	11.77	3.4	ng/ul	180739	2	10.13	1079560	20.0
Benzenesulfonamid...	15.57	4.9	ng/ul	439907	4	16.79	1795990	20.0
Benzenesulfonamid...	16.09	3.1	ng/ul	274732	4	16.79	1795990	20.0
(DEL) Alkane: Str...	20.76	3.4	ng/ul	360816	5	21.01	2144850	20.0