LSC Area Percent Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM022621\
Data File : BM028909.D
Acq On : 25 Feb 2021 22:42
Operator : CG/JU
Sample : M1482-02
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : DBGJ2

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-BM022121MA.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	3.146	24	27	33	rBV	3283	3740	0.23%	0.064%
2	3.904	153	156	160	rBB	1534	2048	0.13%	0.035%
3	5.398	407	410	417	rBB	1888	2803	0.17%	0.048%
4	5.775	470	474	478	rBB2	1417	1746	0.11%	0.030%
5	6.981	675	679	688	rBB2	12688	21661	1.33%	0.368%
6	7.139	700	706	714	rBB	44987	70025	4.29%	1.191%
7	7.328	730	738	748	rBB	54391	86189	5.28%	1.466%
8	7.433	753	756	761	rBB	1955	2526	0.15%	0.043%
9	7.516	765	770	775	rBB	7618	11619	0.71%	0.198%
10	7.798	812	818	825	rBB	57199	82878	5.07%	1.410%
11	8.163	875	880	886	rBB	14889	21999	1.35%	0.374%
12	8.333	903	909	916	rBB	89267	140407	8.59%	2.388%
13	8.528	937	942	950	rBB	37777	56935	3.48%	0.968%
14	8.657	958	964	970	rBB3	5415	9143	0.56%	0.156%
15	8.975	1012	1018	1027	rBB	60643	95696	5.86%	1.628%
16	9.157	1045	1049	1053	rBB2	1562	2302	0.14%	0.039%
17	9.280	1064	1070	1076	rVB2	5690	11392	0.70%	0.194%
18	9.345	1077	1081	1085	rBB2	1724	2315	0.14%	0.039%
19	9.692	1135	1140	1150	rBB	52132	80246	4.91%	1.365%
20	9.792	1153	1157	1161	rBV	4425	6676	0.41%	0.114%
21	9.833	1161	1164	1169	rVB3	1978	3387	0.21%	0.058%
22	9.992	1187	1191	1193	rBV2	2282	3436	0.21%	0.058%
23	10.104	1205	1210	1218	rVB2	4650	9361	0.57%	0.159%
24	10.233	1227	1232	1242	rBB	72334	114431	7.00%	1.946%
25	10.604	1288	1295	1298	rBV	68629	114595	7.01%	1.949%
26	10.657	1298	1304	1315	rVV	961415	1633864	100.00%	27.789%
27	10.774	1317	1324	1331	rVB	52039	87490	5.35%	1.488%
28	10.974	1353	1358	1363	rBB3	2591	3985	0.24%	0.068%
29	11.445	1434	1438	1443	rBB2	1960	2839	0.17%	0.048%
30	11.698	1475	1481	1488	rBB3	1226	2911	0.18%	0.050%
31	12.004	1528	1533	1539	rVB	20817	32923	2.02%	0.560%
32	12.163	1556	1560	1564	rBV	2063	3446	0.21%	0.059%
33	12.268	1572	1578	1585	rBB	127590	190244	11.64%	3.236%
34	12.386	1594	1598	1603	rBV	2144	3592	0.22%	0.061%

LSC Area Percent Report

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Acq On : 25 Feb 2021 22:42
Operator : CG/JU
Sample : M1482-02
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : DBGJ2

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-BM022121MA.M
Title : SVOA CALIBRATION

35	12.486	1606	1615	1622	rBB	72594	114893	7.03%	1.954%
36	12.933	1687	1691	1695	rBV3	1078	2357	0.14%	0.040%
37	13.286	1746	1751	1760	rBB	18058	26770	1.64%	0.455%
38	13.480	1781	1784	1788	rBV	2144	2944	0.18%	0.050%
39	13.615	1804	1807	1809	rBV	2169	2759	0.17%	0.047%
40	13.774	1829	1834	1839	rBB3	4344	6937	0.42%	0.118%
41	13.833	1840	1844	1846	rBV	2434	3253	0.20%	0.055%
42	13.874	1846	1851	1857	rVB	133421	170753	10.45%	2.904%
43	14.151	1892	1898	1908	rBV	150358	213001	13.04%	3.623%
44	14.457	1942	1950	1956	rBV	95856	138024	8.45%	2.348%
45	14.521	1956	1961	1967	rVB	126635	176669	10.81%	3.005%
46	14.604	1970	1975	1981	rBB	20501	27730	1.70%	0.472%
47	14.662	1981	1985	1989	rBB	2324	3130	0.19%	0.053%
48	14.709	1989	1993	1994	rBV2	5399	6673	0.41%	0.113%
49	14.733	1994	1997	2001	rVV	12989	18459	1.13%	0.314%
50	14.768	2001	2003	2007	rVB	2821	3203	0.20%	0.054%
51	14.857	2013	2018	2022	rBV	32245	47051	2.88%	0.800%
52	14.892	2022	2024	2031	rVB	11406	13232	0.81%	0.225%
53	15.015	2042	2045	2048	rBB	1809	2000	0.12%	0.034%
54	15.415	2108	2113	2114	rBV2	2371	2628	0.16%	0.045%
55	15.451	2114	2119	2125	rVV	181345	253936	15.54%	4.319%
56	15.509	2125	2129	2134	rVB	40889	53578	3.28%	0.911%
57	15.598	2140	2144	2153	rBV	58485	77825	4.76%	1.324%
58	15.756	2169	2171	2176	rVB2	1398	1722	0.11%	0.029%
59	15.951	2201	2204	2210	rBB	1652	2147	0.13%	0.037%
60	16.186	2240	2244	2248	rBB2	4536	6591	0.40%	0.112%
61	16.404	2277	2281	2285	rVV	3785	5005	0.31%	0.085%
62	16.480	2289	2294	2301	rVB2	7495	10020	0.61%	0.170%
63	16.833	2350	2354	2357	rBV	3486	4564	0.28%	0.078%
64	16.862	2357	2359	2363	rVB	1650	1817	0.11%	0.031%
65	16.992	2378	2381	2385	rBV	1616	2134	0.13%	0.036%
66	17.168	2407	2411	2412	rBV	1805	2011	0.12%	0.034%
67	17.203	2412	2417	2421	rVV	103543	144219	8.83%	2.453%
68	17.245	2421	2424	2427	rVV	16377	20915	1.28%	0.356%
69	17.298	2427	2433	2440	rVV2	194193	271893	16.64%	4.624%
70	17.351	2440	2442	2448	rVB2	1189	1655	0.10%	0.028%
71	17.574	2476	2480	2482	rBV2	1331	1698	0.10%	0.029%

LSC Area Percent Report

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Misc :
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Instrument : BNA_M
ClientSampleId : DBGJ2

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
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Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM022121MA.M
Title : SVOA CALIBRATION

72	17.609	2482	2486	2494	rVV	95430	129526	7.93%	2.203%
73	17.709	2496	2503	2509	rVV3	10610	16132	0.99%	0.274%
74	17.774	2509	2514	2521	rVV	10603	13776	0.84%	0.234%
75	18.045	2556	2560	2563	rBV	5136	6594	0.40%	0.112%
76	18.086	2563	2567	2570	rBV	11688	12683	0.78%	0.216%
77	18.121	2570	2573	2577	rVB2	4094	5230	0.32%	0.089%
78	18.203	2582	2587	2594	rVV	6758	9759	0.60%	0.166%
79	18.362	2609	2614	2616	rBV3	1778	2708	0.17%	0.046%
80	18.386	2616	2618	2622	rVV3	1571	2398	0.15%	0.041%
81	18.421	2622	2624	2628	rVV	1709	2282	0.14%	0.039%
82	18.521	2637	2641	2647	rBV3	2298	4114	0.25%	0.070%
83	18.580	2647	2651	2656	rBV2	1455	1977	0.12%	0.034%
84	19.221	2755	2760	2763	rBV2	1506	2082	0.13%	0.035%
85	19.362	2781	2784	2787	rBV3	3546	3669	0.22%	0.062%
86	19.586	2816	2822	2828	rVB	221999	281697	17.24%	4.791%
87	19.633	2828	2830	2834	rBV2	1438	1944	0.12%	0.033%
88	20.003	2890	2893	2896	rVB2	1804	2217	0.14%	0.038%
89	20.062	2901	2903	2909	rVB3	2507	3780	0.23%	0.064%
90	20.603	2993	2995	2998	rBV4	1549	1731	0.11%	0.029%
91	21.068	3070	3074	3079	rBV	27213	32102	1.96%	0.546%
92	21.362	3119	3124	3129	rBV	118952	142247	8.71%	2.419%
93	23.432	3470	3476	3484	rVB	162780	285384	17.47%	4.854%
94	23.568	3494	3499	3509	rVB2	83707	152472	9.33%	2.593%

Sum of corrected areas: 5879550

LSC Report - Integrated Chromatogram

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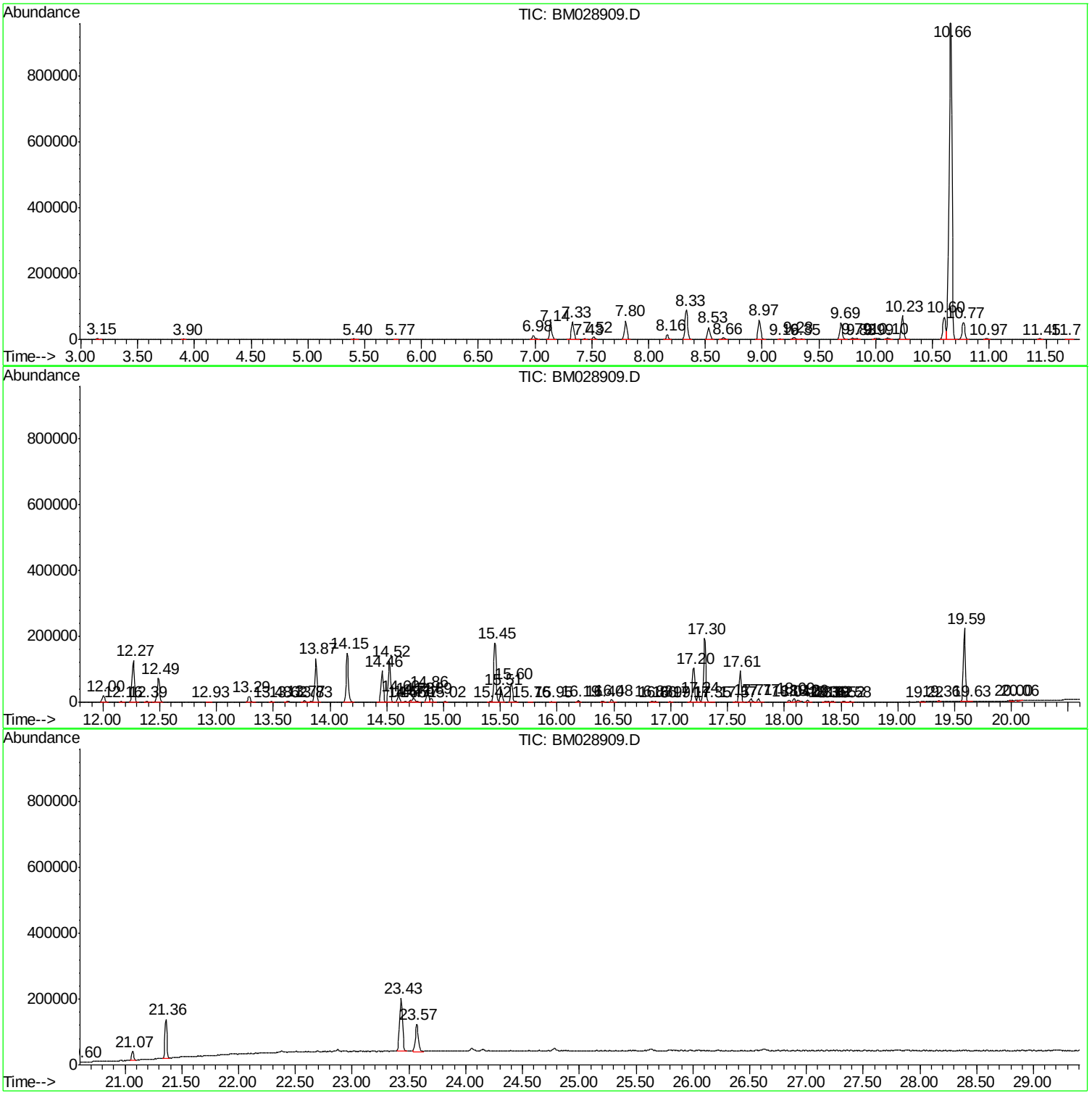
Data File : BM028909.D
Acq On : 25 Feb 2021 22:42
Operator : CG/JU
Sample : M1482-02
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : DBGJ2

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM022121MA.M
Quant Title : SVOA CALIBRATION

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



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM022621\
Data File : BM028909.D
Acq On : 25 Feb 2021 22:42
Operator : CG/JU
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Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : DBGJ2

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM022121MA.M
Quant Title : SVOA CALIBRATION

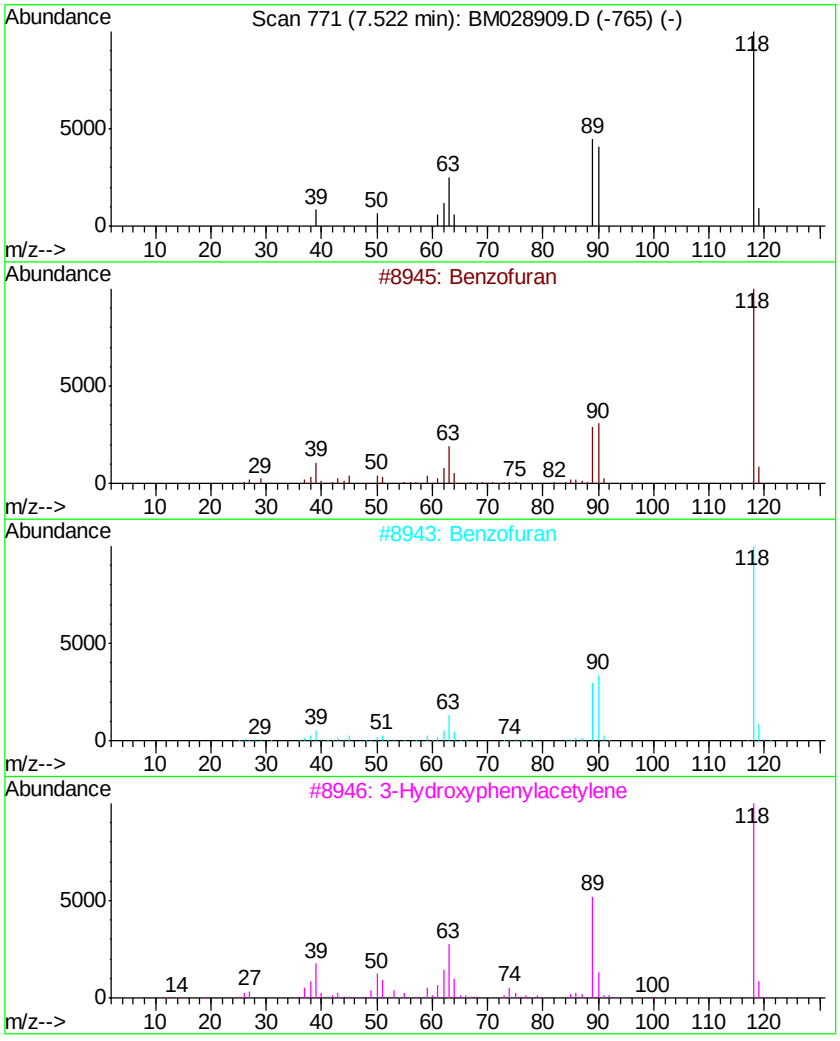
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

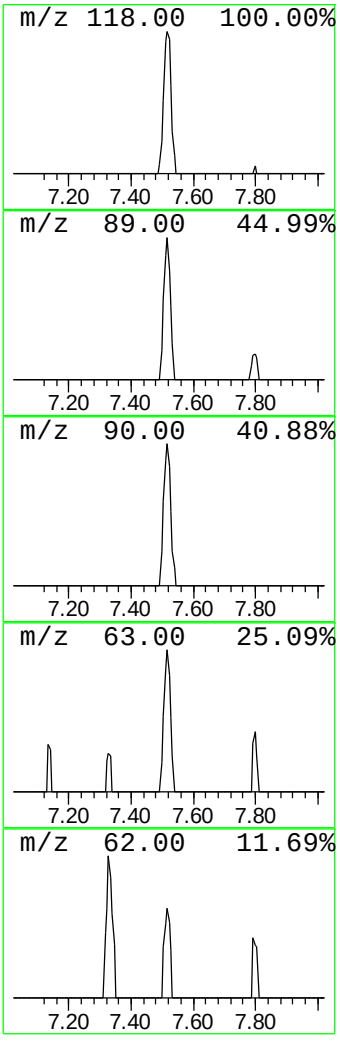
****

Peak Number 1 Benzofuran Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.52	2.80 ng/ul	11619	1,4-Dichlorobenzene-d4	7.80

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzofuran	118	C8H6O	000271-89-6	86
2	Benzofuran	118	C8H6O	000271-89-6	72
3	3-Hydroxyphenylacetylene	118	C8H6O	010401-11-3	64
4	1H-Pyrrolo[2,3-b]pyridine	118	C7H6N2	000271-63-6	53
5	3-Pyridineacetonitrile	118	C7H6N2	006443-85-2	50





Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM022621\
Data File : BM028909.D
Acq On : 25 Feb 2021 22:42
Operator : CG/JU
Sample : M1482-02
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : DBGJ2

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM022121MA.M
Quant Title : SVOA CALIBRATION

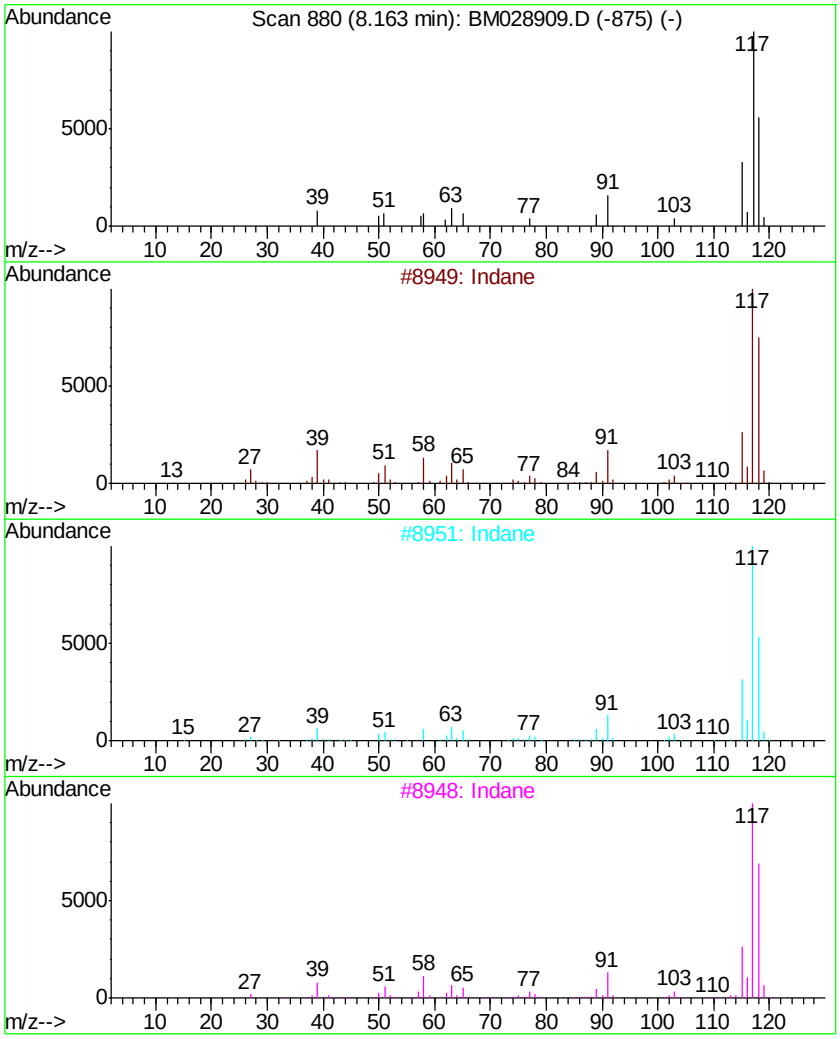
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

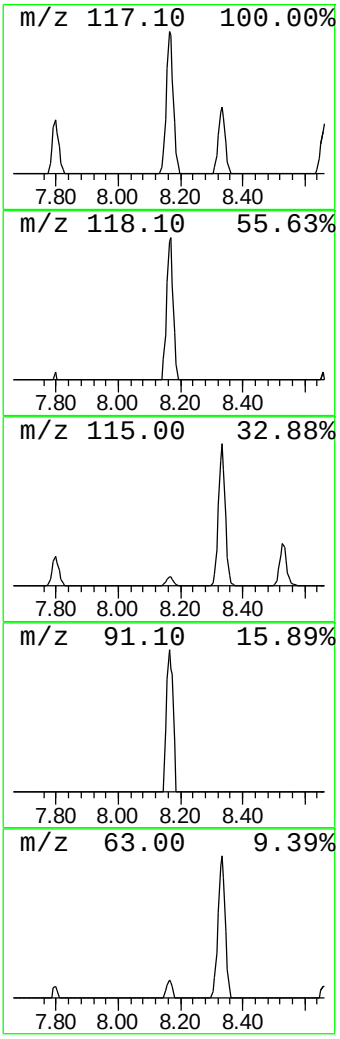
****

Peak Number 2 Indane Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.16	5.31 ng/ul	21999	1,4-Dichlorobenzene-d4	7.80

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indane	118	C9H10	000496-11-7	90
2	Indane	118	C9H10	000496-11-7	87
3	Indane	118	C9H10	000496-11-7	72
4	Tetracyclo[3.3.1.0(2,8).0(4,6)]-...	118	C9H10	1000191-13-7	72
5	Benzene, cyclopropyl-	118	C9H10	000873-49-4	64





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Instrument : BNA_M
ClientSampleId : DBGJ2

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM022121MA.M
Quant Title : SVOA CALIBRATION

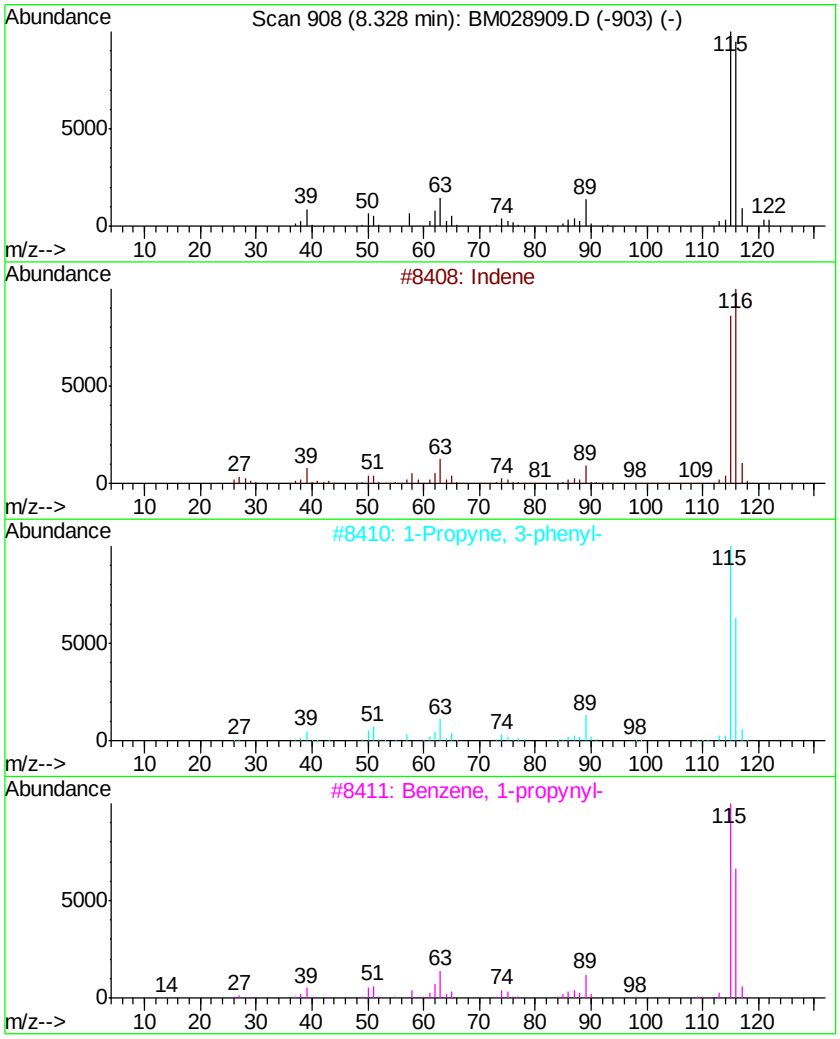
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

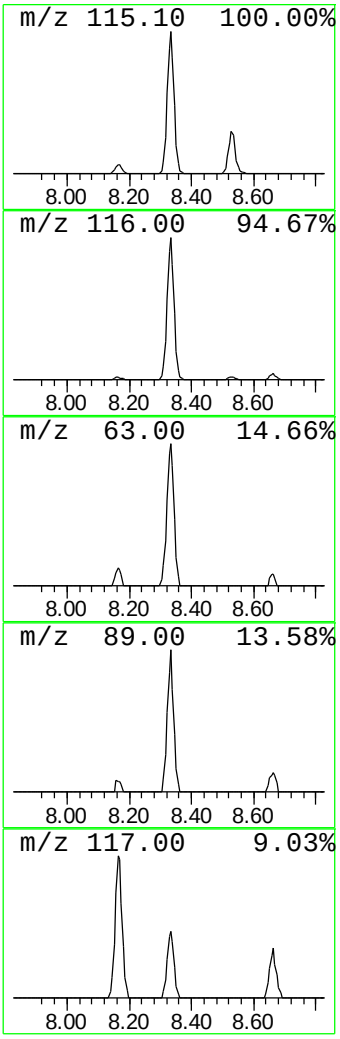
****

Peak Number 3 1-Propyne, 3-phenyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.33	33.88 ng/ul	140407	1,4-Dichlorobenzene-d4	7.80

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indene	116	C9H8	000095-13-6	95
2	1-Propyne, 3-phenyl-	116	C9H8	010147-11-2	95
3	Benzene, 1-propynyl-	116	C9H8	000673-32-5	94
4	Benzene, 1,2-propadienyl-	116	C9H8	002327-99-3	91
5	3-Methylphenylacetylene	116	C9H8	000766-82-5	91





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Instrument : BNA_M
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM022121MA.M
Quant Title : SVOA CALIBRATION

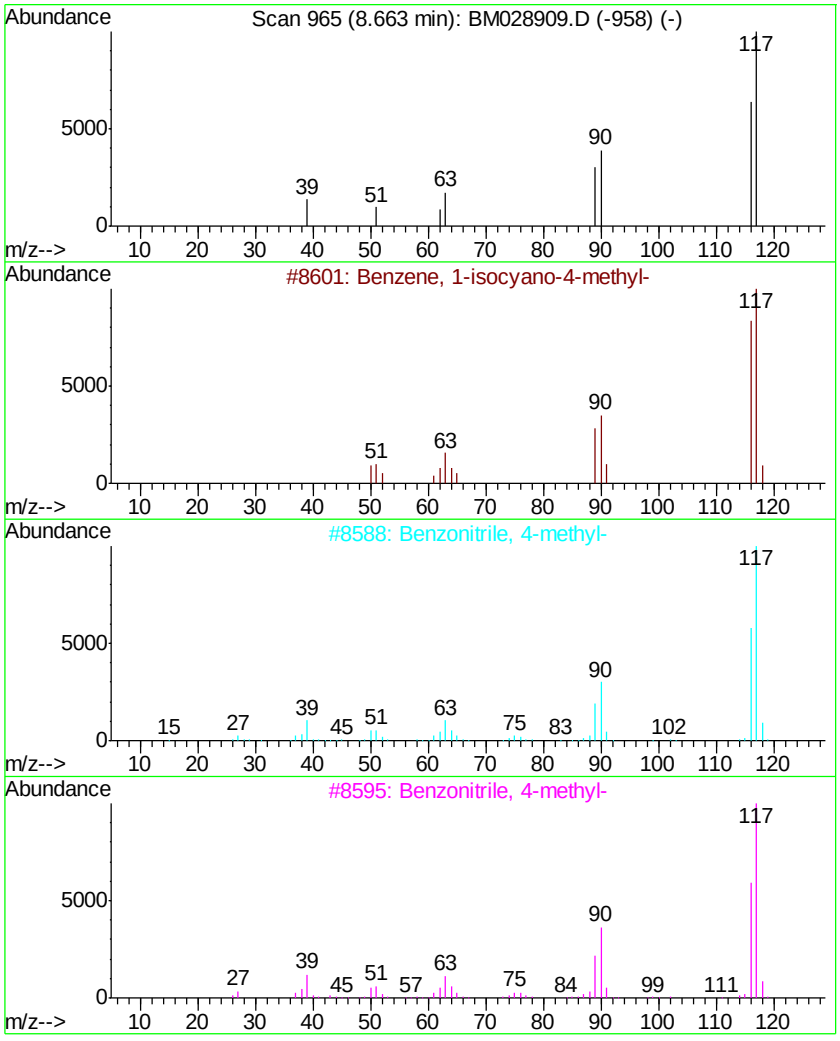
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

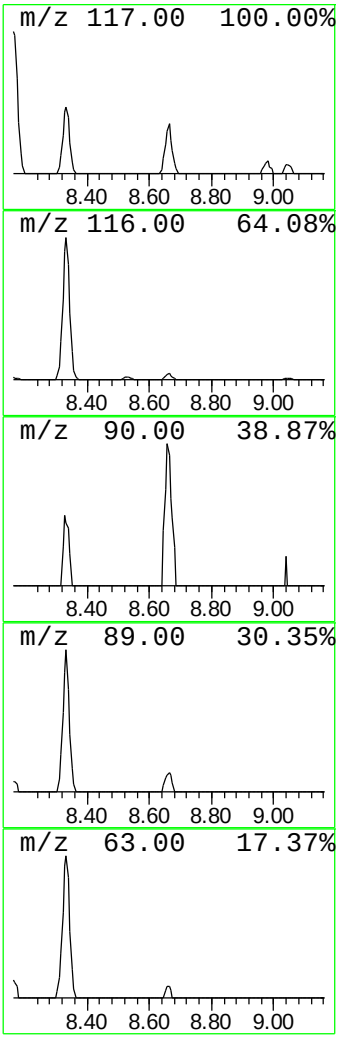
****

Peak Number 4 Benzene, 1-isocyano-4-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.66	2.21 ng/ul	9143	1,4-Dichlorobenzene-d4	7.80

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-isocyano-4-methyl-	117	C8H7N	007175-47-5	91
2	Benzonitrile, 4-methyl-	117	C8H7N	000104-85-8	91
3	Benzonitrile, 4-methyl-	117	C8H7N	000104-85-8	91
4	Benzonitrile, 4-methyl-	117	C8H7N	000104-85-8	91
5	Benzyl nitrile	117	C8H7N	000140-29-4	91





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Sample : M1482-02
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : DBGJ2

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM022121MA.M
Quant Title : SVOA CALIBRATION

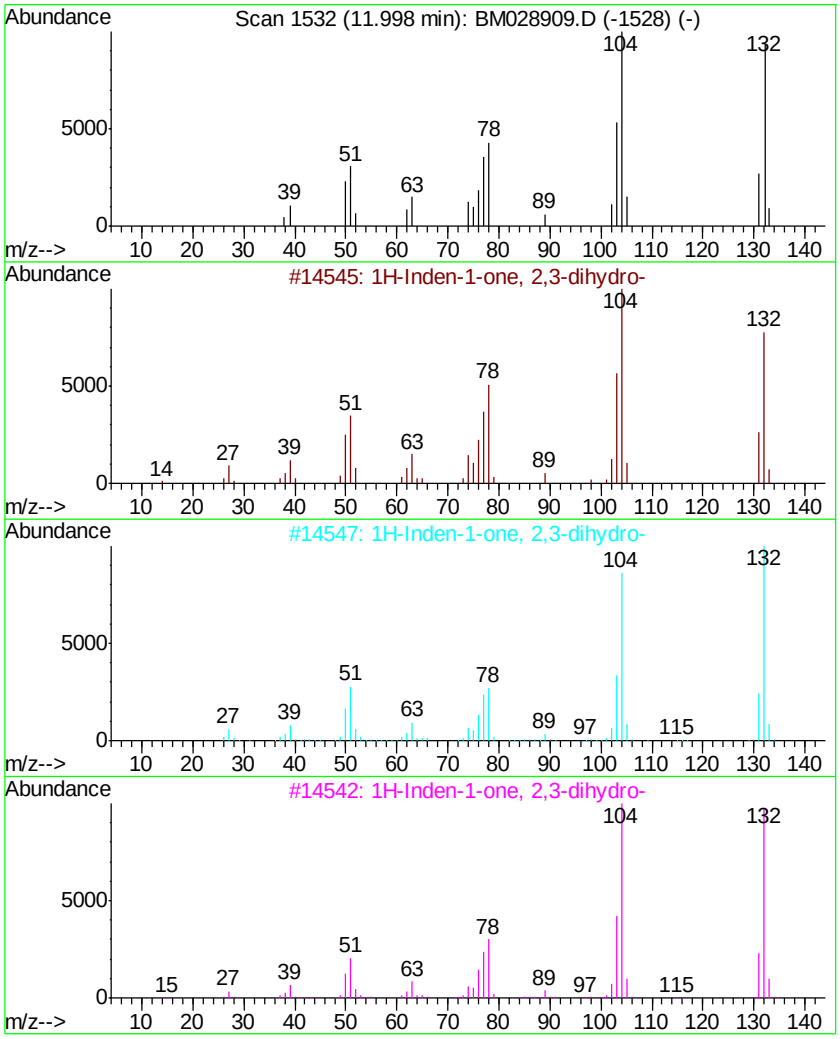
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

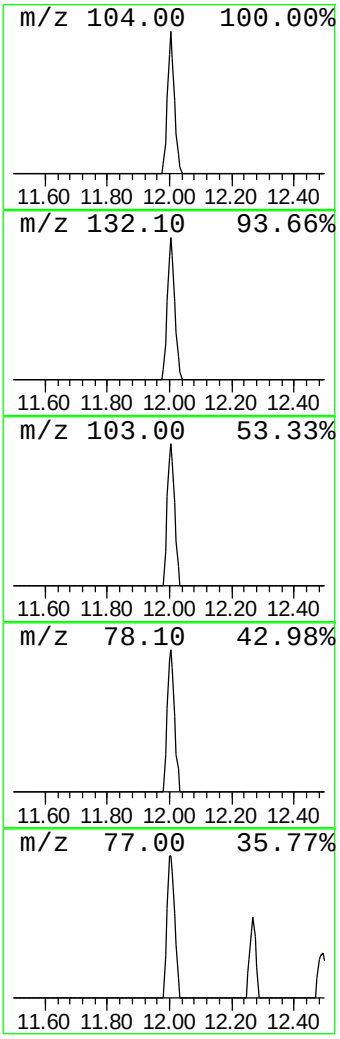
****

Peak Number 5 1H-Inden-1-one, 2,3-dihydro- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.00	5.75 ng/ul	32923	Naphthalene-d8	10.60

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Inden-1-one, 2,3-dihydro-	132	C9H8O	000083-33-0	96
2	1H-Inden-1-one, 2,3-dihydro-	132	C9H8O	000083-33-0	95
3	1H-Inden-1-one, 2,3-dihydro-	132	C9H8O	000083-33-0	95
4	2H-Inden-2-one, 1,3-dihydro-	132	C9H8O	000615-13-4	64
5	2H-Inden-2-one, 1,3-dihydro-	132	C9H8O	000615-13-4	64





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Instrument : BNA_M
ClientSampleId : DBGJ2

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM022121MA.M
Quant Title : SVOA CALIBRATION

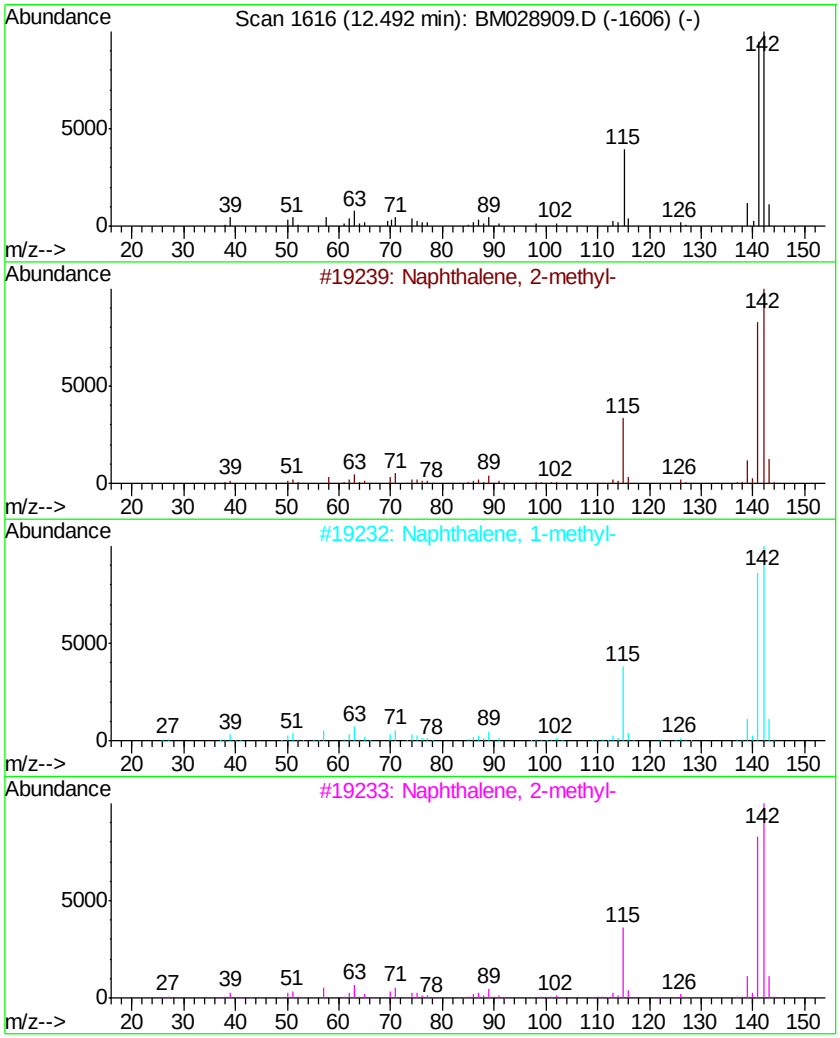
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

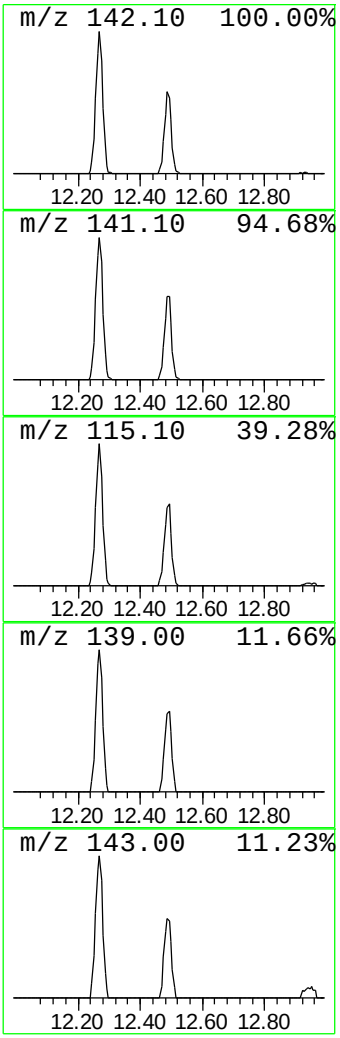
****

Peak Number 6 Naphthalene, 1-methyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.49	20.05 ng/ul	114893	Naphthalene-d8	10.60

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 2-methyl-	142	C11H10	000091-57-6	97
2	Naphthalene, 1-methyl-	142	C11H10	000090-12-0	96
3	Naphthalene, 2-methyl-	142	C11H10	000091-57-6	96
4	Naphthalene, 2-methyl-	142	C11H10	000091-57-6	94
5	Naphthalene, 1-methyl-	142	C11H10	000090-12-0	94





Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM022621\
Data File : BM028909.D
Acq On : 25 Feb 2021 22:42
Operator : CG/JU
Sample : M1482-02
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : DBGJ2

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM022121MA.M
Quant Title : SVOA CALIBRATION

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

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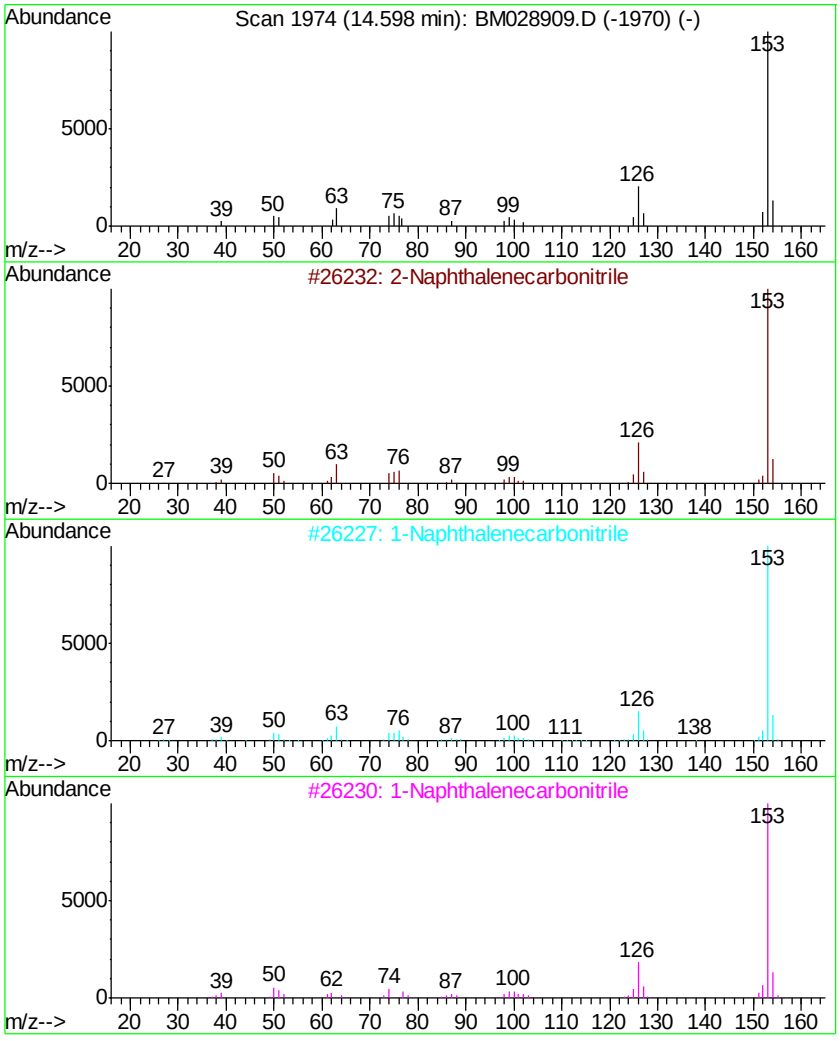
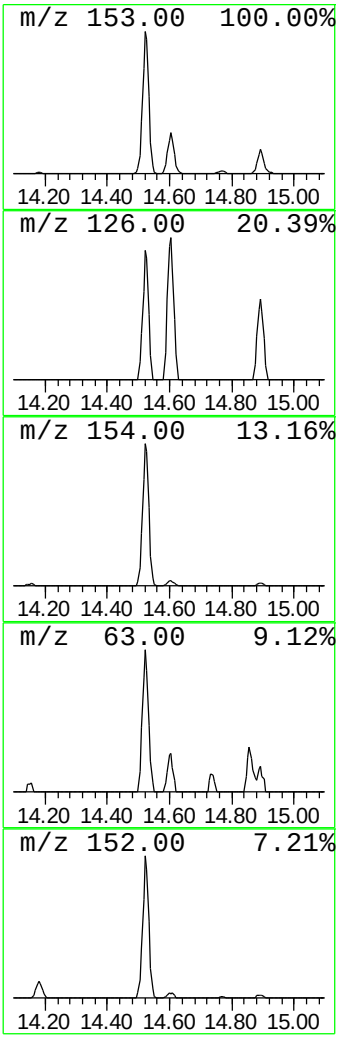
Peak Number 7 2-Naphthalenecarbonitrile Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.60	4.02 ng/ul	27730	Acenaphthene-d10	14.46

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Naphthalenecarbonitrile	153	C11H7N	000613-46-7	95
2	1-Naphthalenecarbonitrile	153	C11H7N	000086-53-3	95
3	1-Naphthalenecarbonitrile	153	C11H7N	000086-53-3	91
4	1-Naphthalenecarbonitrile	153	C11H7N	000086-53-3	91
5	2-Naphthalenecarbonitrile	153	C11H7N	000613-46-7	91





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Operator : CG/JU
Sample : M1482-02
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : DBGJ2

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM022121MA.M
Quant Title : SVOA CALIBRATION

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

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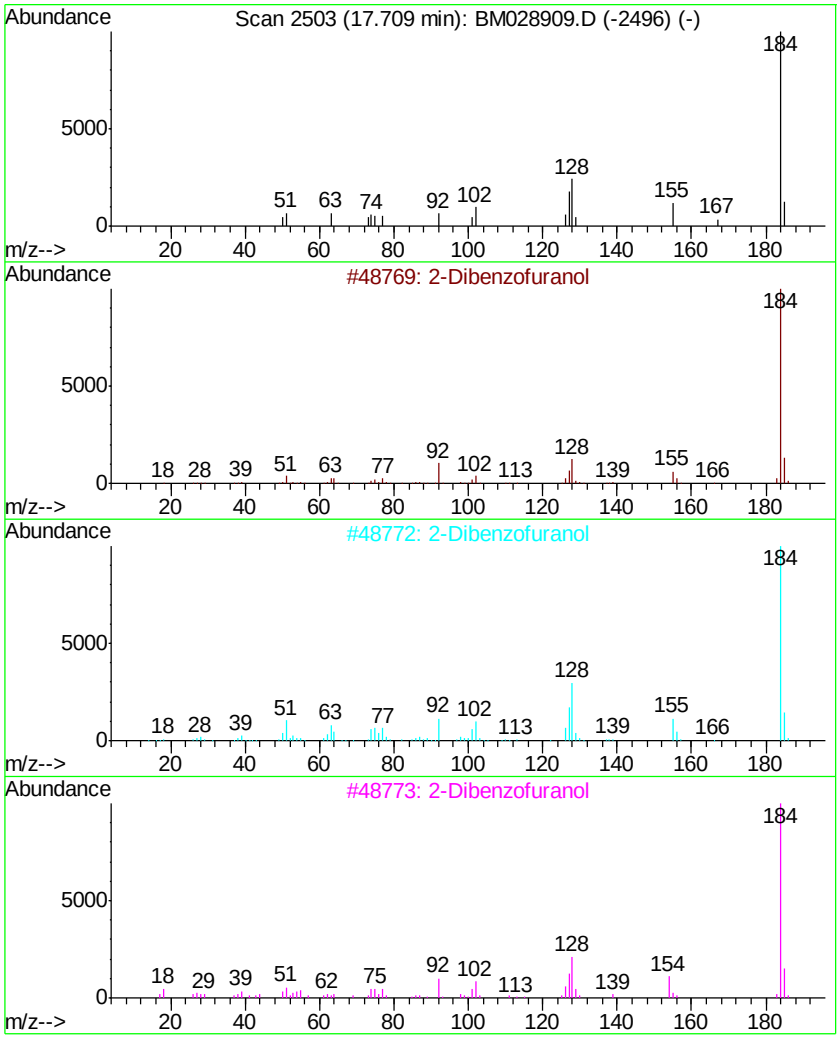
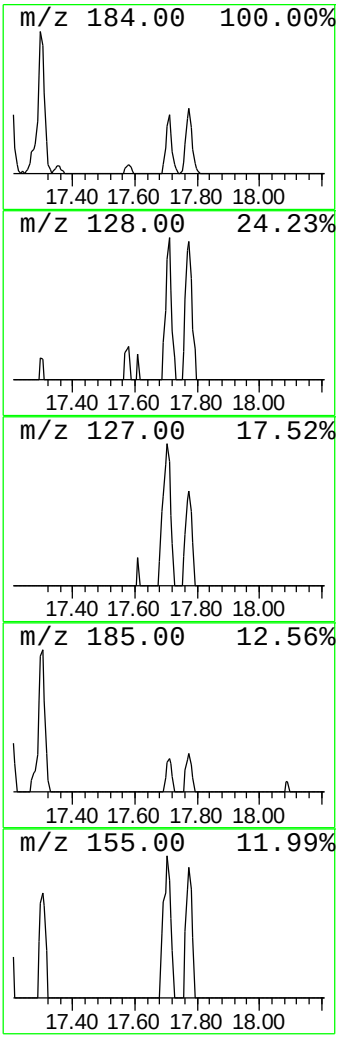
Peak Number 8 2-Dibenzofuranol Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.71	2.24 ng/ul	16132	Phenanthrene-d10	17.20

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Dibenzofuranol	184	C12H8O2	000086-77-1	91
2	2-Dibenzofuranol	184	C12H8O2	000086-77-1	91
3	2-Dibenzofuranol	184	C12H8O2	000086-77-1	86
4	Dibenzo-p-dioxin	184	C12H8O2	000262-12-4	78
5	1H-Pyrazolo[3,4-b]quinolin-3-amine	184	C10H8N4	1000319-54-1	59





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Data File : BM028909.D
Acq On : 25 Feb 2021 22:42
Operator : CG/JU
Sample : M1482-02
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : DBGJ2

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM022121MA.M
Quant Title : SVOA CALIBRATION

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

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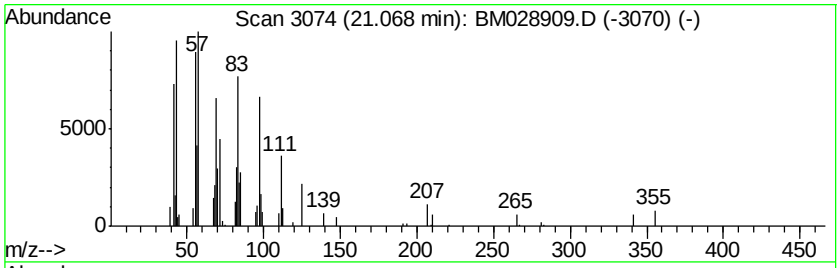
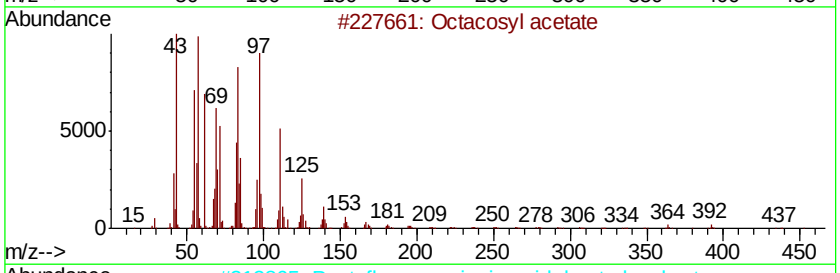
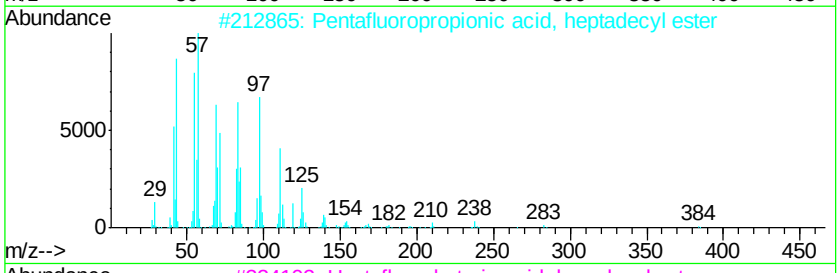
Peak Number 9 Octacosyl acetate Concentration Rank 5

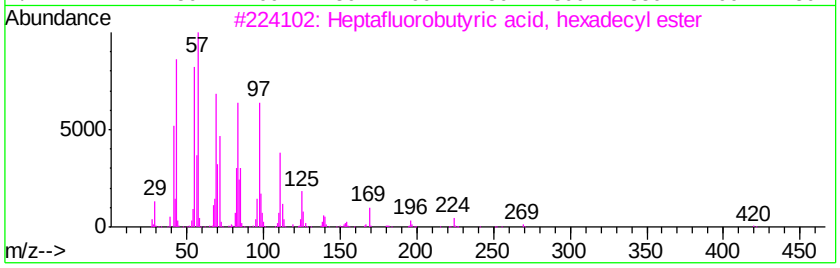
R.T.	EstConc	Area	Relative to ISTD	R.T.
21.07	4.51 ng/ul	32102	Chrysene-d12	21.36

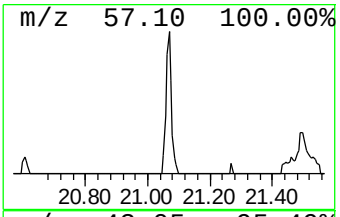
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octacosyl acetate	452	C30H60O2	018206-97-8	93
2	Pentafluoropropionic acid, hepta...	402	C20H35F5O2	959218-78-1	91
3	Heptafluorobutyric acid, hexadec...	438	C20H33F7O2	006385-15-5	91
4	1-Tricosene	322	C23H46	018835-32-0	90
5	Pentafluoropropionic acid, penta...	374	C18H31F5O2	959092-08-1	90

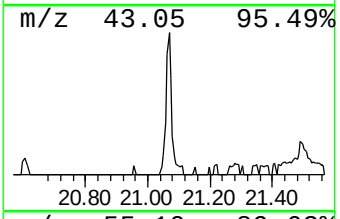


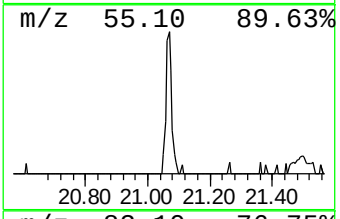


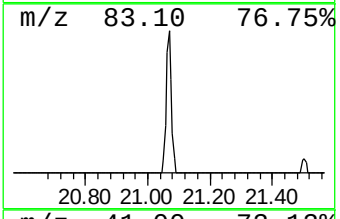


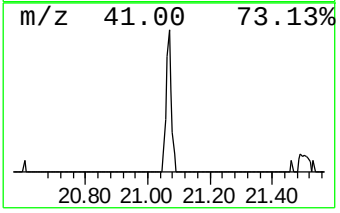












Tentatively Identified Compound (LSC) summary

Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM022621\
Data File : BM028909.D
Acq On : 25 Feb 2021 22:42
Operator : CG/JU
Sample : M1482-02
Misc :
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Instrument : BNA_M
ClientSampleId : DBGJ2

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM022121MA.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
Benzofuran	7.52	2.8	ng/ul	11619	1	7.80	82878	20.0
Indane	8.16	5.3	ng/ul	21999	1	7.80	82878	20.0
1-Propyne, 3-phenyl-	8.33	33.9	ng/ul	140407	1	7.80	82878	20.0
Benzene, 1-isocya...	8.66	2.2	ng/ul	9143	1	7.80	82878	20.0
1H-Inden-1-one, 2...	12.00	5.8	ng/ul	32923	2	10.60	114595	20.0
Naphthalene, 1-me...	12.49	20.1	ng/ul	114893	2	10.60	114595	20.0
2-Naphthalenecarb...	14.60	4.0	ng/ul	27730	3	14.46	138024	20.0
2-Dibenzofuranol	17.71	2.2	ng/ul	16132	4	17.20	144219	20.0
Octacosyl acetate	21.07	4.5	ng/ul	32102	5	21.36	142247	20.0