LSC Area Percent Report

Data Path : \\74.0.250.170\SVOASRV\HPCHEM1\BNA M\DATA\BM121417\
Data File : BM013406.D
Acq On : 14 Dec 2017 17:34
Operator : SJ/JU
Sample : I6900-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : F9C00

Integration Parameters: LSCINT.P
Integrator: RTE
Smoothing : OFF Filtering: 5
Sampling : 1 Min Area: 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:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM113017.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	5.752	448	453	459	rVB2	144033	218948	3.36%	0.344%
2	6.857	632	641	649	rBV	143338	251424	3.85%	0.395%
3	7.022	662	669	676	rBV	565705	894244	13.71%	1.406%
4	7.210	695	701	709	rBV	599863	993595	15.23%	1.562%
5	7.675	773	780	787	rBV	645461	1035494	15.87%	1.628%
6	7.746	787	792	798	rVB2	52352	88497	1.36%	0.139%
7	8.187	861	867	874	rBV2	55047	90613	1.39%	0.142%
8	8.381	894	900	910	rBV	464844	770547	11.81%	1.212%
9	8.834	969	977	985	rBV	829510	1366373	20.94%	2.149%
10	9.093	1015	1021	1029	rBV2	141435	220929	3.39%	0.347%
11	9.545	1092	1098	1106	rBV	713616	1217317	18.66%	1.914%
12	9.640	1109	1114	1124	rBV	186466	340167	5.21%	0.535%
13	9.745	1126	1132	1137	rBV4	40328	74843	1.15%	0.118%
14	9.828	1137	1146	1150	rBV3	40200	83060	1.27%	0.131%
15	9.951	1160	1167	1171	rBV	801335	1369029	20.98%	2.153%
16	9.992	1171	1174	1182	rVV4	213027	418166	6.41%	0.658%
17	10.081	1182	1189	1197	rVV2	1042115	2073394	31.78%	3.260%
18	10.140	1197	1199	1205	rVB4	65964	96383	1.48%	0.152%
19	10.457	1245	1253	1260	rBV	908895	1469941	22.53%	2.311%
20	11.257	1381	1389	1392	rBV2	160972	360070	5.52%	0.566%
21	11.292	1392	1395	1402	rVB	191359	291949	4.47%	0.459%
22	11.716	1461	1467	1473	rBV4	77367	129923	1.99%	0.204%
23	11.986	1507	1513	1520	rBV5	40397	82500	1.26%	0.130%
24	12.228	1549	1554	1560	rBV	166103	265844	4.07%	0.418%
25	12.381	1576	1580	1586	rVV3	72096	135087	2.07%	0.212%
26	12.457	1586	1593	1599	rVB7	55933	117715	1.80%	0.185%
27	13.063	1689	1696	1702	rBV9	28839	72173	1.11%	0.113%
28	13.739	1804	1811	1817	rBV	2095185	2971339	45.54%	4.672%
29	13.845	1823	1829	1837	rBV	578133	843259	12.92%	1.326%
30	14.010	1847	1857	1868	rBV	2349908	3530504	54.11%	5.551%
31	14.104	1868	1873	1882	rVB3	36566	79677	1.22%	0.125%
32	14.322	1901	1910	1917	rBV2	1387747	2178752	33.39%	3.426%
33	14.380	1917	1920	1927	rVB2	99869	145368	2.23%	0.229%
34	14.539	1941	1947	1953	rBV	161658	281720	4.32%	0.443%

LSC Area Percent Report

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Acq On : 14 Dec 2017 17:34
Operator : SJ/JU
Sample : I6900-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : F9C00

Integration Parameters: LSCINT.P
Integrator: RTE
Smoothing : OFF Filtering: 5
Sampling : 1 Min Area: 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:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM113017.M
Title : SVOA CALIBRATION

35	14.598	1953	1957	1967	rVB4	148494	280953	4.31%	0.442%
36	14.927	2006	2013	2023	rBV	146802	267675	4.10%	0.421%
37	15.251	2061	2068	2074	rBV	319967	522694	8.01%	0.822%
38	15.316	2074	2079	2086	rVV	2959320	4675378	71.66%	7.352%
39	15.374	2086	2089	2095	rVB3	94608	137132	2.10%	0.216%
40	15.445	2095	2101	2107	rBV	1171968	1642959	25.18%	2.583%
41	15.551	2115	2119	2124	rVB3	44285	71155	1.09%	0.112%
42	16.045	2188	2203	2211	rBV5	357927	791957	12.14%	1.245%
43	16.121	2211	2216	2220	rVV2	112399	203143	3.11%	0.319%
44	16.163	2220	2223	2226	rVV2	80141	124600	1.91%	0.196%
45	16.210	2226	2231	2236	rVV4	110631	234286	3.59%	0.368%
46	16.263	2236	2240	2248	rVB3	182385	324303	4.97%	0.510%
47	16.339	2248	2253	2257	rBV	99038	131899	2.02%	0.207%
48	16.727	2315	2319	2323	rVB	60304	91475	1.40%	0.144%
49	16.857	2336	2341	2345	rBV2	57192	95046	1.46%	0.149%
50	17.068	2368	2377	2382	rBV2	2097530	3050266	46.75%	4.796%
51	17.110	2382	2384	2387	rVV	118783	142476	2.18%	0.224%
52	17.168	2387	2394	2400	rVB2	3513956	4931195	75.58%	7.754%
53	17.445	2435	2441	2443	rBV3	68178	103538	1.59%	0.163%
54	17.474	2443	2446	2453	rVV	80086	119858	1.84%	0.188%
55	17.568	2455	2462	2468	rBV2	80438	159952	2.45%	0.252%
56	17.633	2468	2473	2483	rVB	83236	141036	2.16%	0.222%
57	17.974	2528	2531	2533	rBV3	49125	66708	1.02%	0.105%
58	18.139	2552	2559	2564	rVB3	41428	75893	1.16%	0.119%
59	18.880	2680	2685	2691	rVB	190161	271884	4.17%	0.428%
60	19.133	2724	2728	2735	rVB	106495	152860	2.34%	0.240%
61	19.257	2741	2749	2756	rVB5	117858	251439	3.85%	0.395%
62	19.351	2761	2765	2770	rVB3	52957	85013	1.30%	0.134%
63	19.468	2779	2785	2791	rBV	4564448	5966894	91.45%	9.382%
64	21.262	3084	3090	3096	rBV	3077206	3807558	58.35%	5.987%
65	23.345	3436	3444	3455	rVB2	3509318	6524829	100.00%	10.260%
66	23.486	3461	3468	3475	rVB	1945496	3591389	55.04%	5.647%

Sum of corrected areas: 63596287

LSC Report - Integrated Chromatogram

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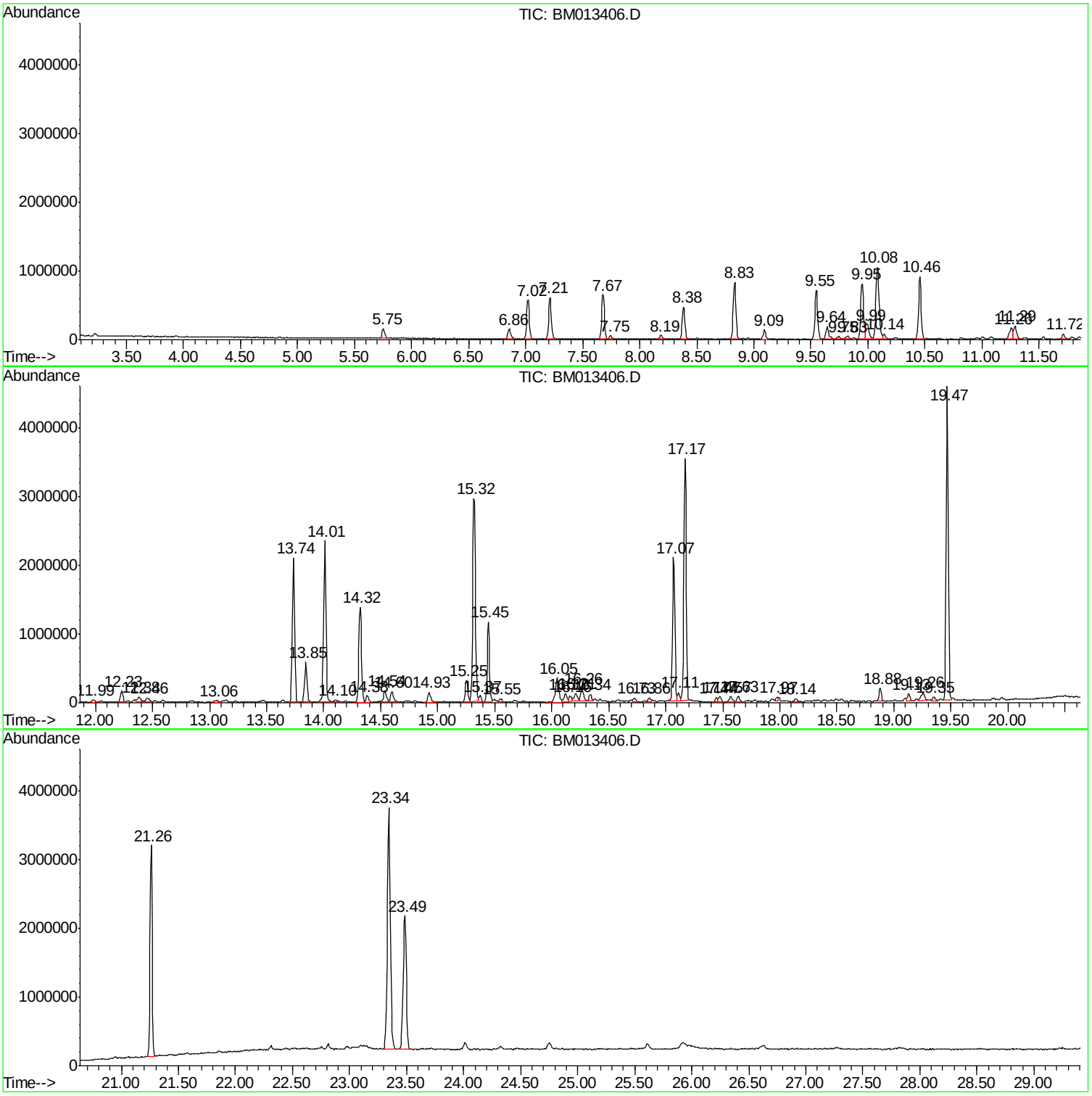
Data File : BM013406.D
Acq On : 14 Dec 2017 17:34
Operator : SJ/JU
Sample : I6900-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : F9C00

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM113017.M
Quant Title : SVOA CALIBRATION

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



Library Search Compound Report

Data Path : \\74.0.250.170\SVOASRV\HPCHEM1\BNA M\DATA\BM121417\
Data File : BM013406.D
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Quant Title : SVOA CALIBRATION

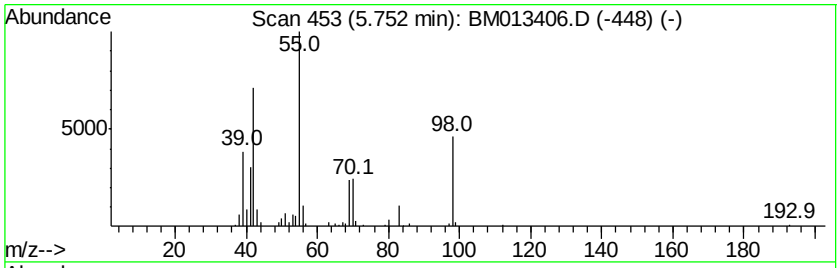
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

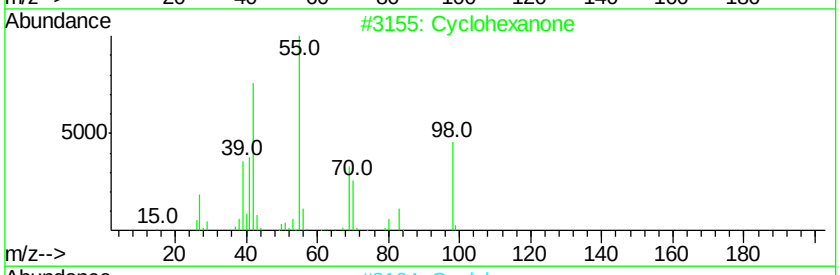
****

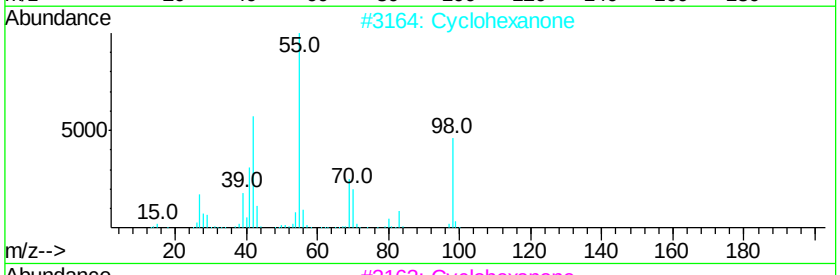
Peak Number 1 Cyclohexanone Concentration Rank 7

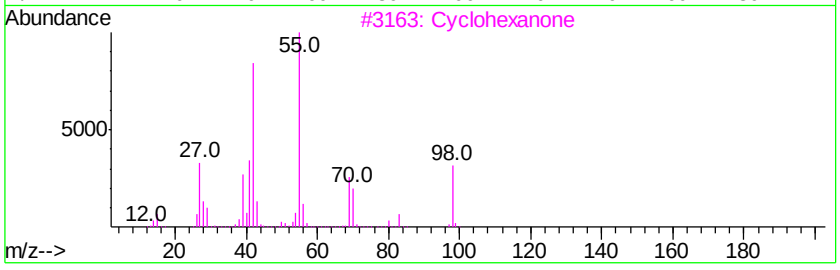
R.T.	EstConc	Area	Relative to ISTD	R.T.
5.75	4.23 ng/ul	218948	1,4-Dichlorobenzene-d4	7.68

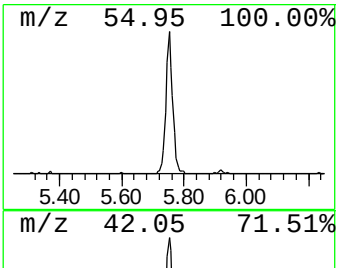
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclohexanone	98	C6H10O	000108-94-1	70
2	Cyclohexanone	98	C6H10O	000108-94-1	62
3	Cyclohexanone	98	C6H10O	000108-94-1	59
4	Cyclohexanone	98	C6H10O	000108-94-1	59
5	Cyclohexanone	98	C6H10O	000108-94-1	56

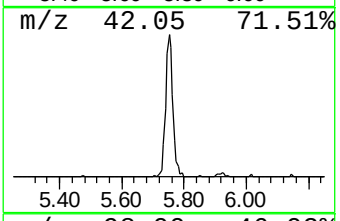


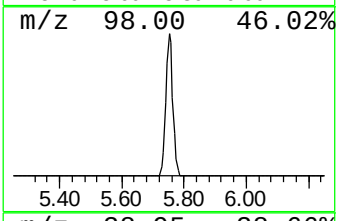


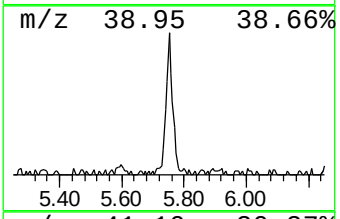


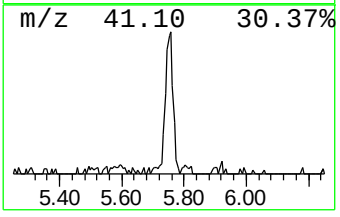












Library Search Compound Report

Data Path : \\74.0.250.170\SVOASRV\HPCHEM1\BNA M\DATA\BM121417\
Data File : BM013406.D
Acq On : 14 Dec 2017 17:34
Operator : SJ/JU
Sample : I6900-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : F9C00

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM113017.M
Quant Title : SVOA CALIBRATION

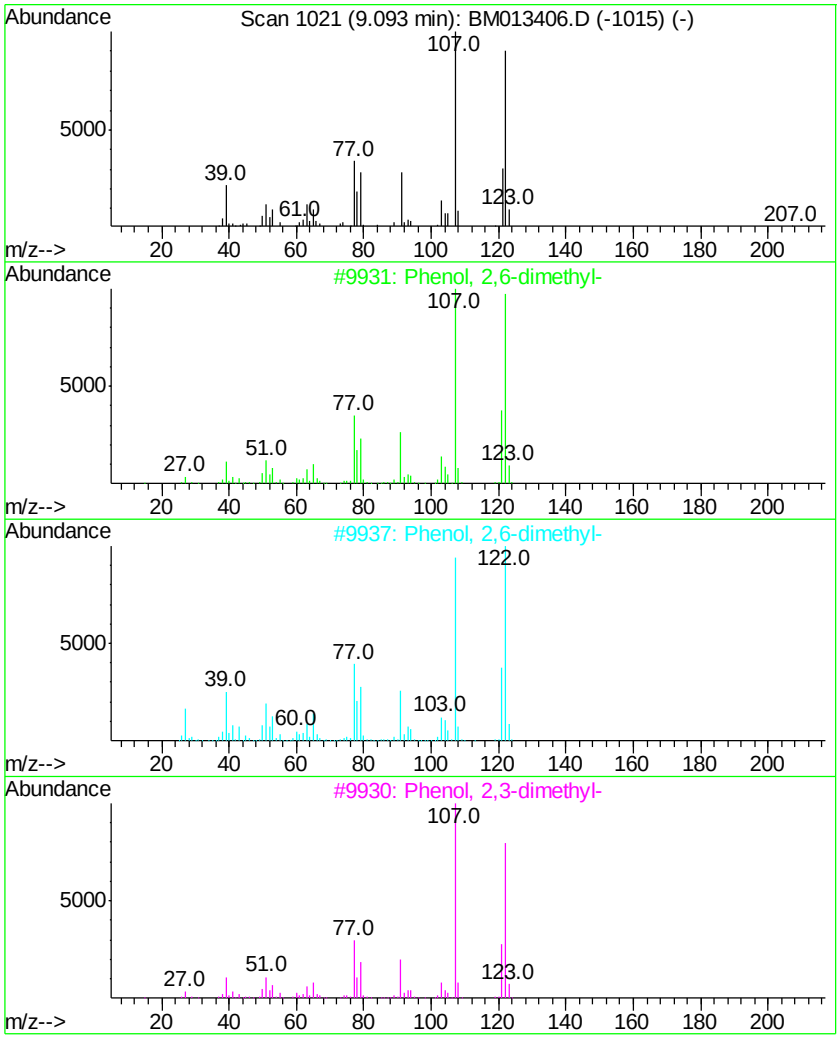
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

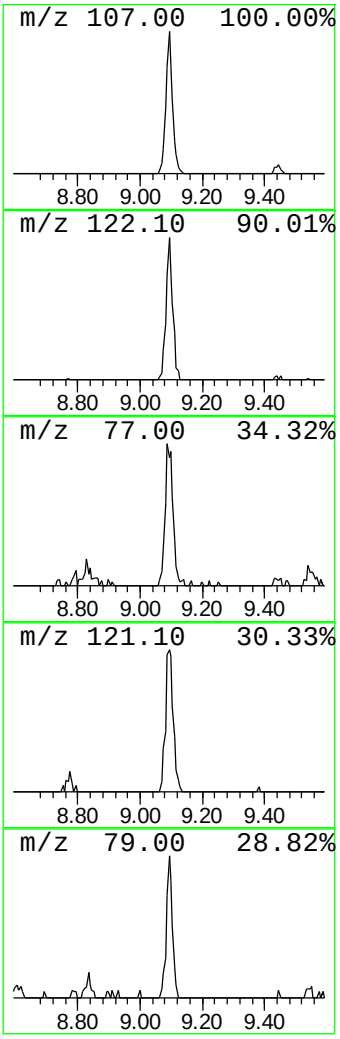
****

Peak Number 2 Phenol, 2,6-dimethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.09	3.01 ng/ul	220929	Naphthalene-d8	10.46

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, 2,6-dimethyl-	122	C8H10O	000576-26-1	96
2	Phenol, 2,6-dimethyl-	122	C8H10O	000576-26-1	95
3	Phenol, 2,3-dimethyl-	122	C8H10O	000526-75-0	94
4	Phenol, 2,3-dimethyl-	122	C8H10O	000526-75-0	94
5	Phenol, 2,6-dimethyl-	122	C8H10O	000576-26-1	93





Library Search Compound Report

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Sample : I6900-01
Misc :
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Instrument : BNA_M
ClientSampleId : F9C00

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM113017.M
Quant Title : SVOA CALIBRATION

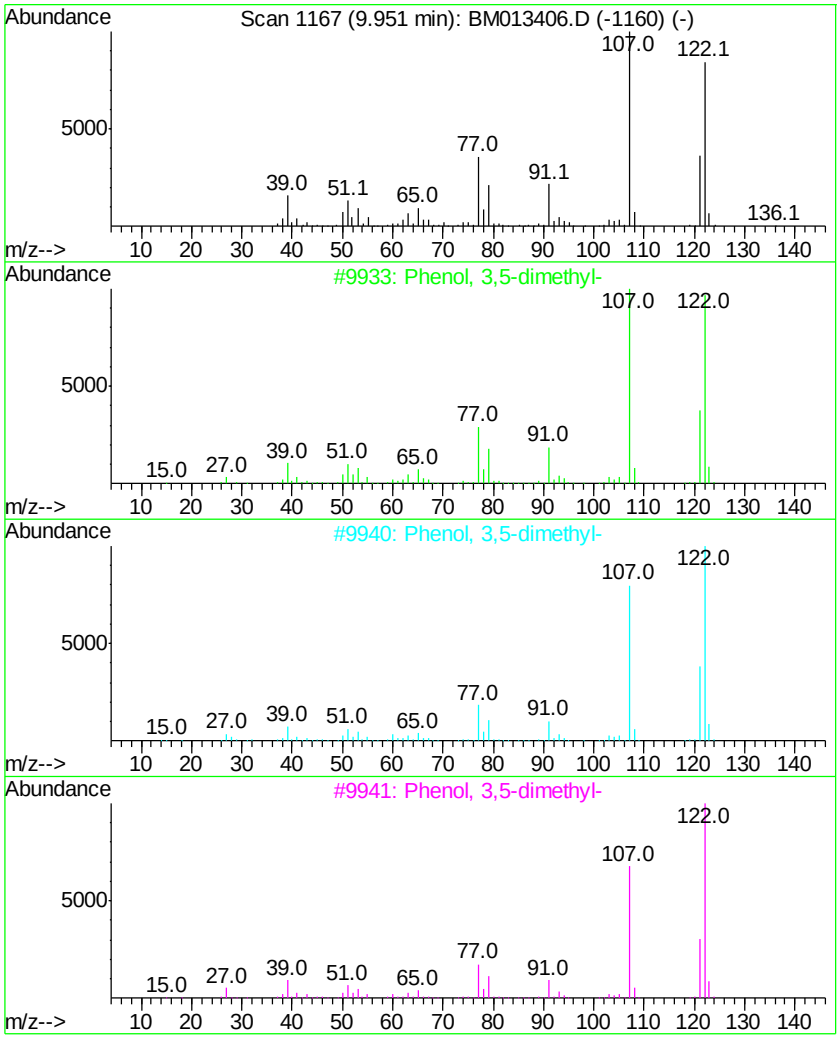
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

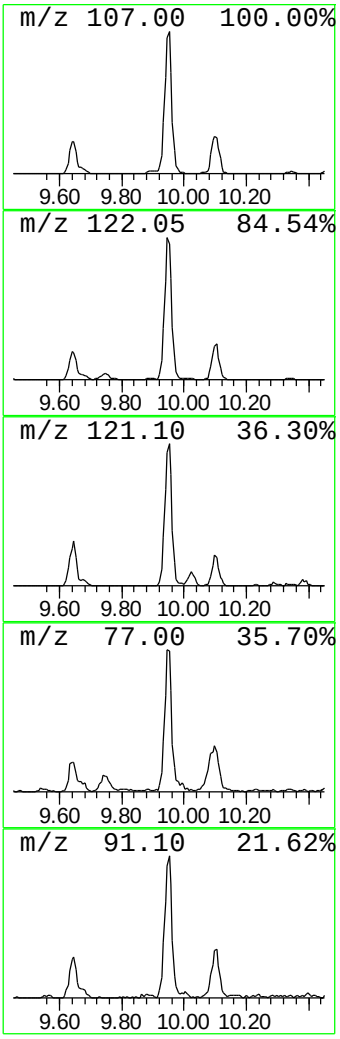
****

Peak Number 3 Phenol, 3,5-dimethyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.95	18.63 ng/ul	1369030	Naphthalene-d8	10.46

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, 3,5-dimethyl-	122	C8H10O	000108-68-9	96
2	Phenol, 3,5-dimethyl-	122	C8H10O	000108-68-9	95
3	Phenol, 3,5-dimethyl-	122	C8H10O	000108-68-9	93
4	Phenol, 3,5-dimethyl-	122	C8H10O	000108-68-9	93
5	Phenol, 2,4-dimethyl-	122	C8H10O	000105-67-9	93





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Sample : I6900-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : F9C00

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM113017.M
Quant Title : SVOA CALIBRATION

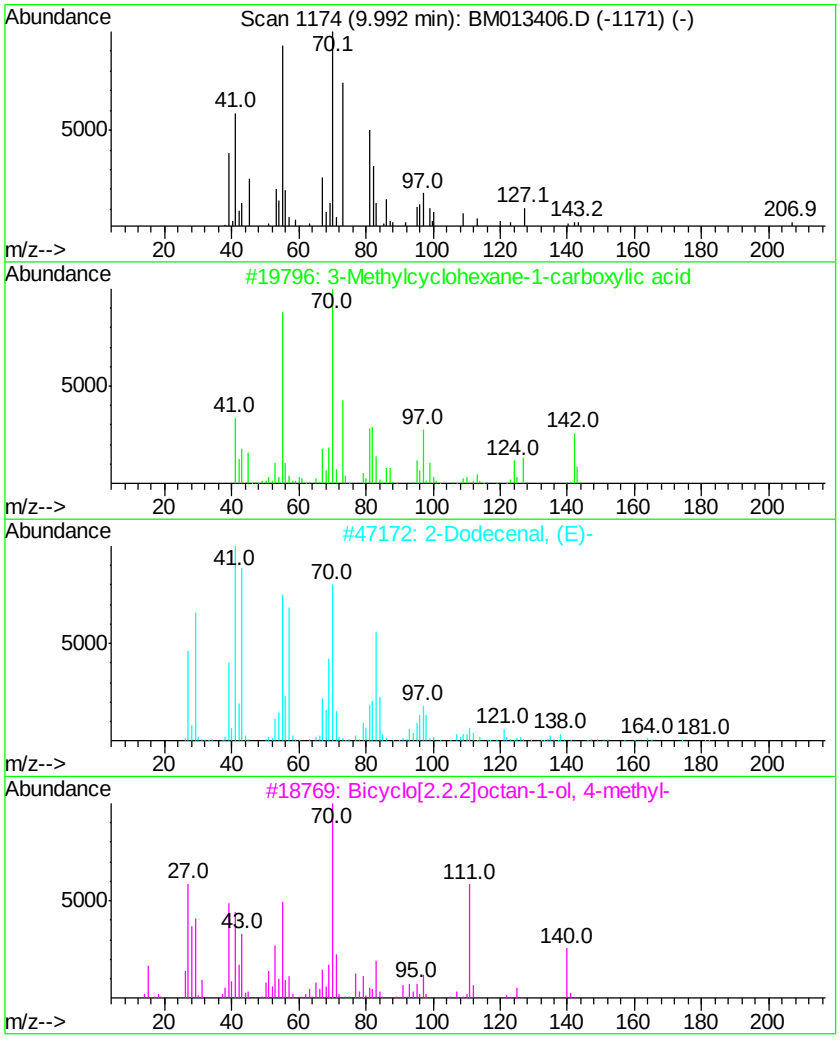
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

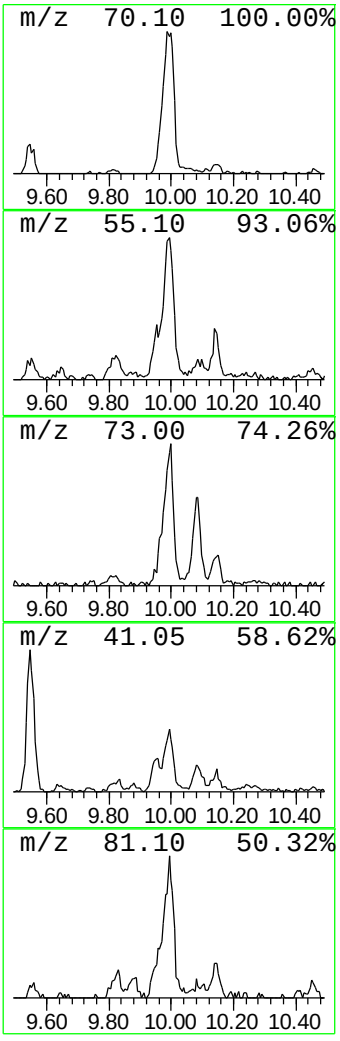
****

Peak Number 4 3-Methylcyclohexane-1-carbo... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.99	5.69 ng/ul	418166	Naphthalene-d8	10.46

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Methylcyclohexane-1-carboxylic...	142	C8H14O2	1000132-15-7	72
2	2-Dodecenal, (E)-	182	C12H22O	020407-84-5	38
3	Bicyclo[2.2.2]octan-1-ol, 4-methyl-	140	C9H16O	000824-13-5	32
4	trans-2-undecenoic acid	184	C11H20O2	015790-94-0	27
5	2-Butene, 2-methyl-	70	C5H10	000513-35-9	27





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Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : F9C00

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM113017.M
Quant Title : SVOA CALIBRATION

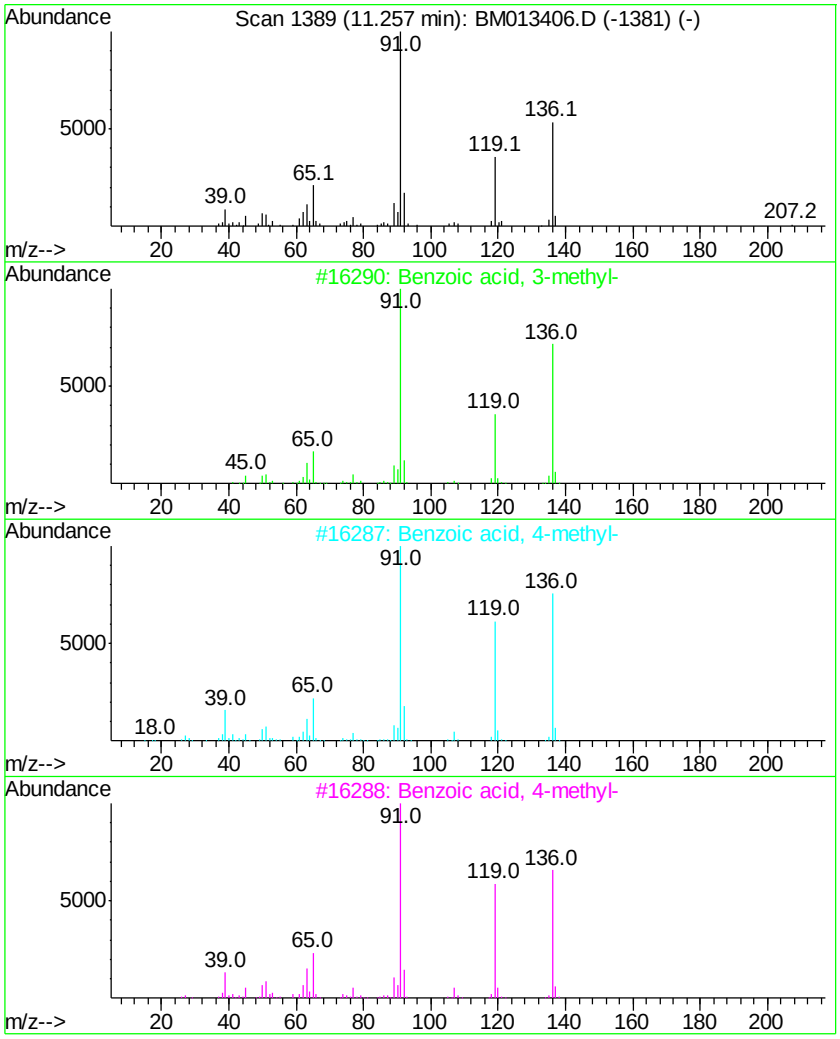
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

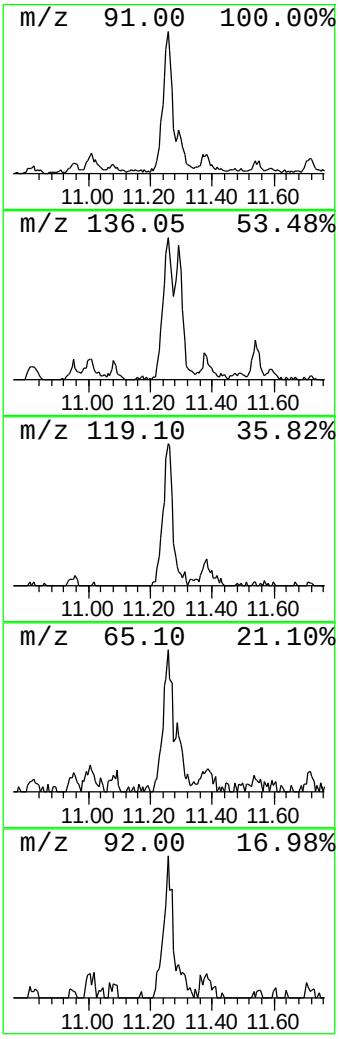
****

Peak Number 5 Benzoic acid, 3-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.26	4.90 ng/ul	360070	Naphthalene-d8	10.46

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzoic acid, 3-methyl-	136	C8H8O2	000099-04-7	95
2	Benzoic acid, 4-methyl-	136	C8H8O2	000099-94-5	91
3	Benzoic acid, 4-methyl-	136	C8H8O2	000099-94-5	87
4	Benzeneacetic acid	136	C8H8O2	000103-82-2	72
5	Benzoic acid, 3-methyl-	136	C8H8O2	000099-04-7	72





Library Search Compound Report

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Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : F9C00

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM113017.M
Quant Title : SVOA CALIBRATION

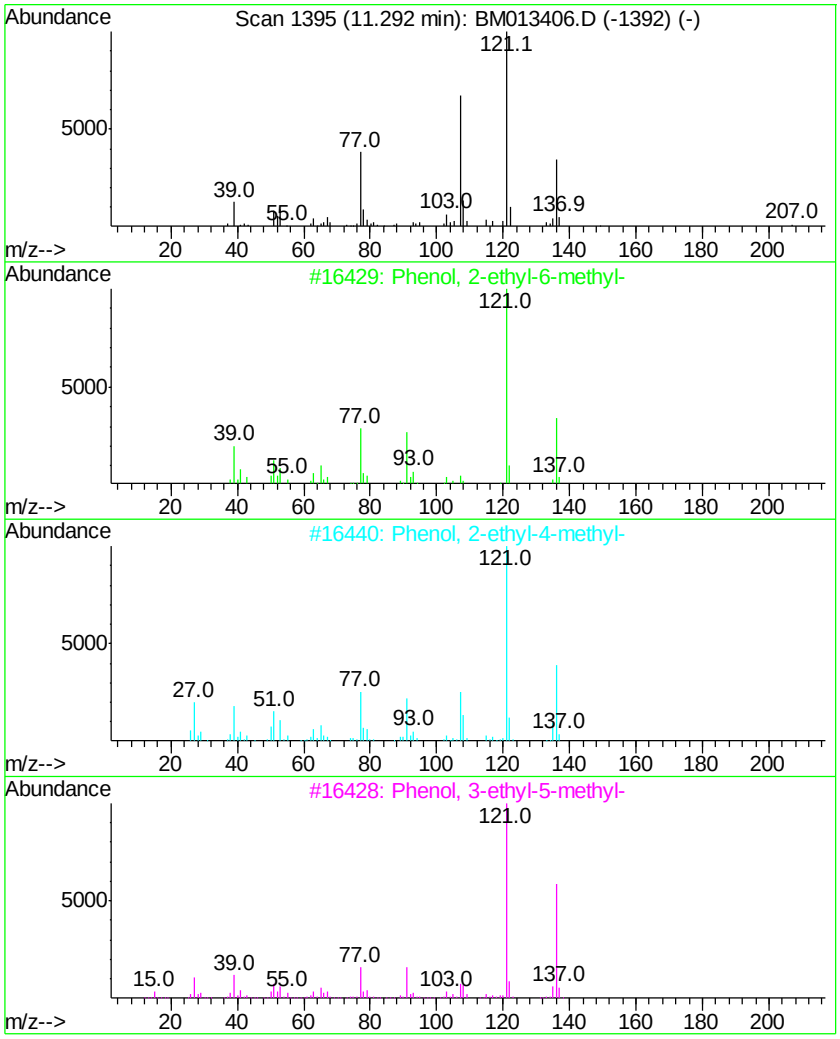
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

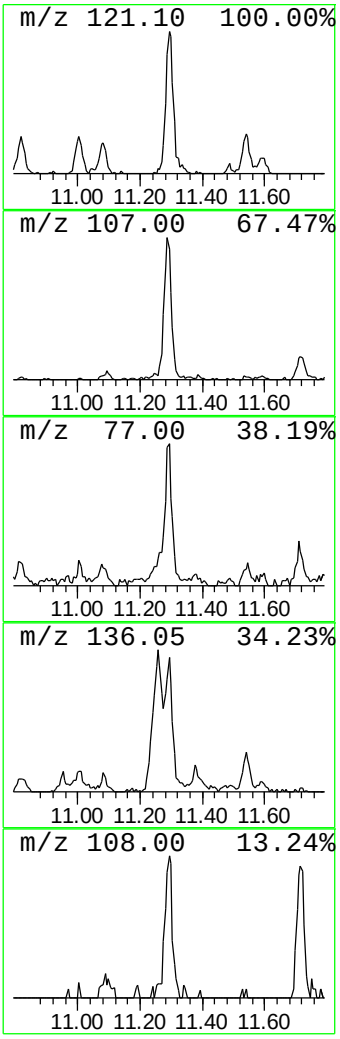
****

Peak Number 6 Phenol, 2-ethyl-6-methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.29	3.97 ng/ul	291949	Naphthalene-d8	10.46

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, 2-ethyl-6-methyl-	136	C9H12O	001687-64-5	72
2	Phenol, 2-ethyl-4-methyl-	136	C9H12O	003855-26-3	68
3	Phenol, 3-ethyl-5-methyl-	136	C9H12O	000698-71-5	59
4	Phenol, 3-(1-methylethyl)-	136	C9H12O	000618-45-1	59
5	Phenol, 4-(1-methylethyl)-	136	C9H12O	000099-89-8	59





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Acq On : 14 Dec 2017 17:34
Operator : SJ/JU
Sample : I6900-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : F9C00

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM113017.M
Quant Title : SVOA CALIBRATION

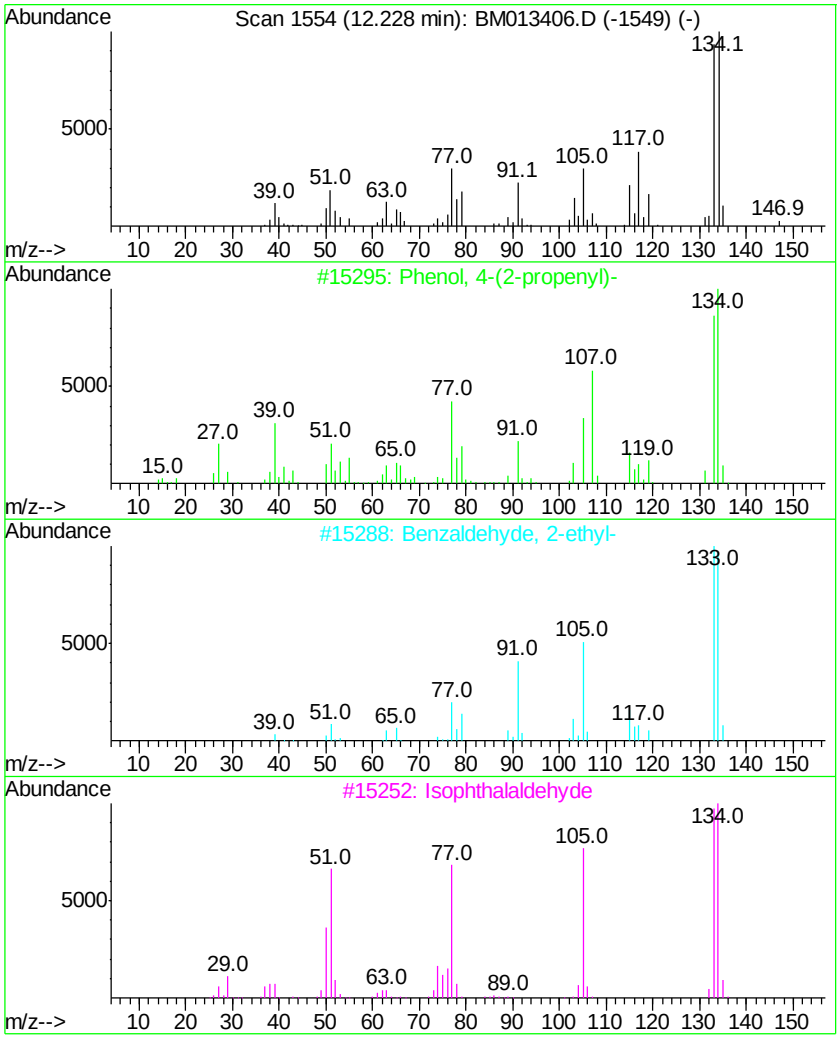
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

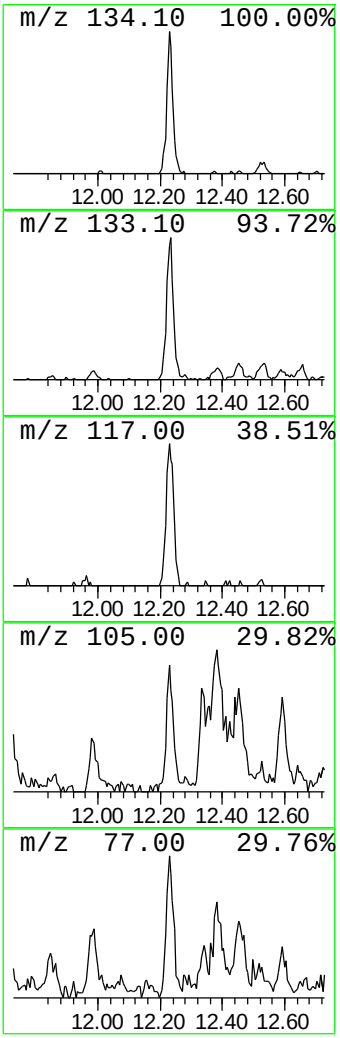
****

Peak Number 7 Phenol, 4-(2-propenyl)- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.23	3.62 ng/ul	265844	Naphthalene-d8	10.46

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, 4-(2-propenyl)-	134	C9H10O	000501-92-8	83
2	Benzaldehyde, 2-ethyl-	134	C9H10O	022927-13-5	74
3	Isophthalaldehyde	134	C8H6O2	000626-19-7	70
4	Phenol, 4-(2-propenyl)-	134	C9H10O	000501-92-8	64
5	Isophthalaldehyde	134	C8H6O2	000626-19-7	64





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Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : F9C00

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM113017.M
Quant Title : SVOA CALIBRATION

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

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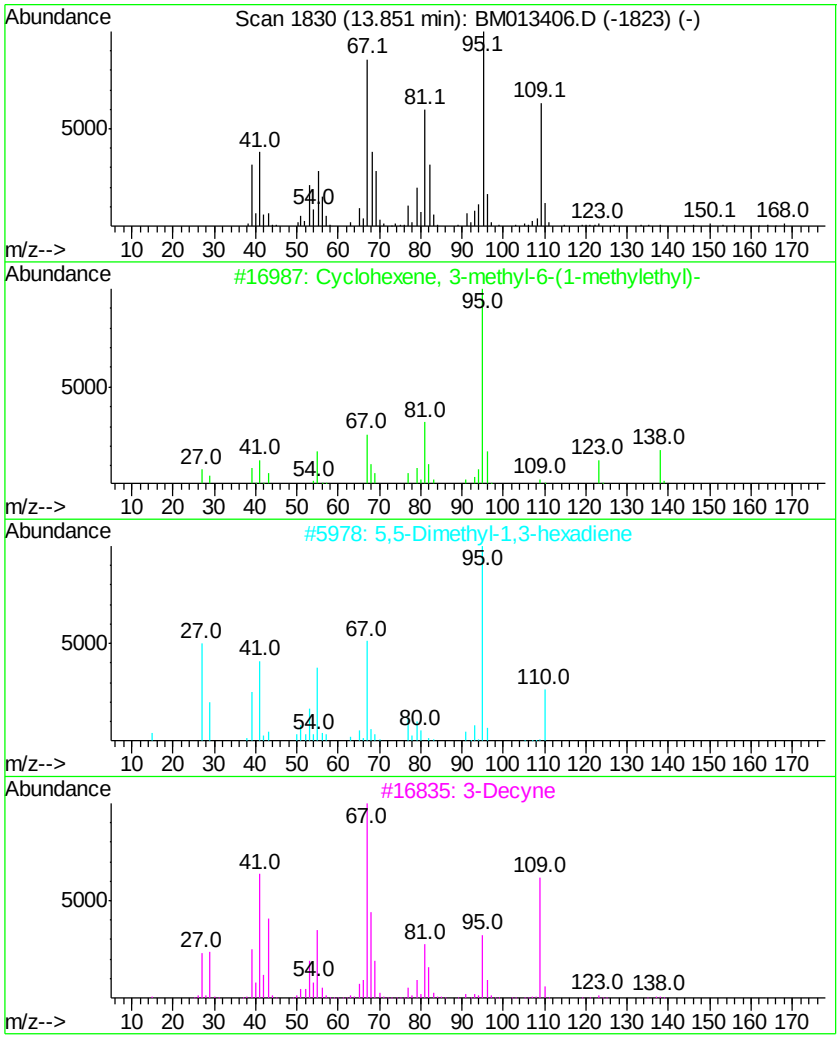
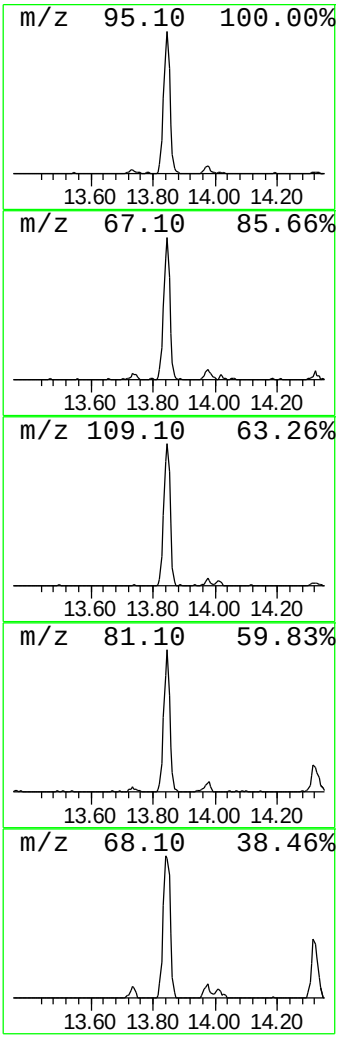
Peak Number 8 Cyclohexene, 3-methyl-6-(1-... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.85	7.74 ng/ul	843259	Acenaphthene-d10	14.32

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclohexene, 3-methyl-6-(1-methy...	138	C10H18	005256-65-5	76
2	5,5-Dimethyl-1,3-hexadiene	110	C8H14	001515-79-3	58
3	3-Decyne	138	C10H18	002384-85-2	58
4	3-Decyne	138	C10H18	002384-85-2	58
5	Bicyclo[3.1.0]hexane, 1,5-dimethyl-	110	C8H14	1000142-17-5	53





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Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : F9C00

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM113017.M
Quant Title : SVOA CALIBRATION

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

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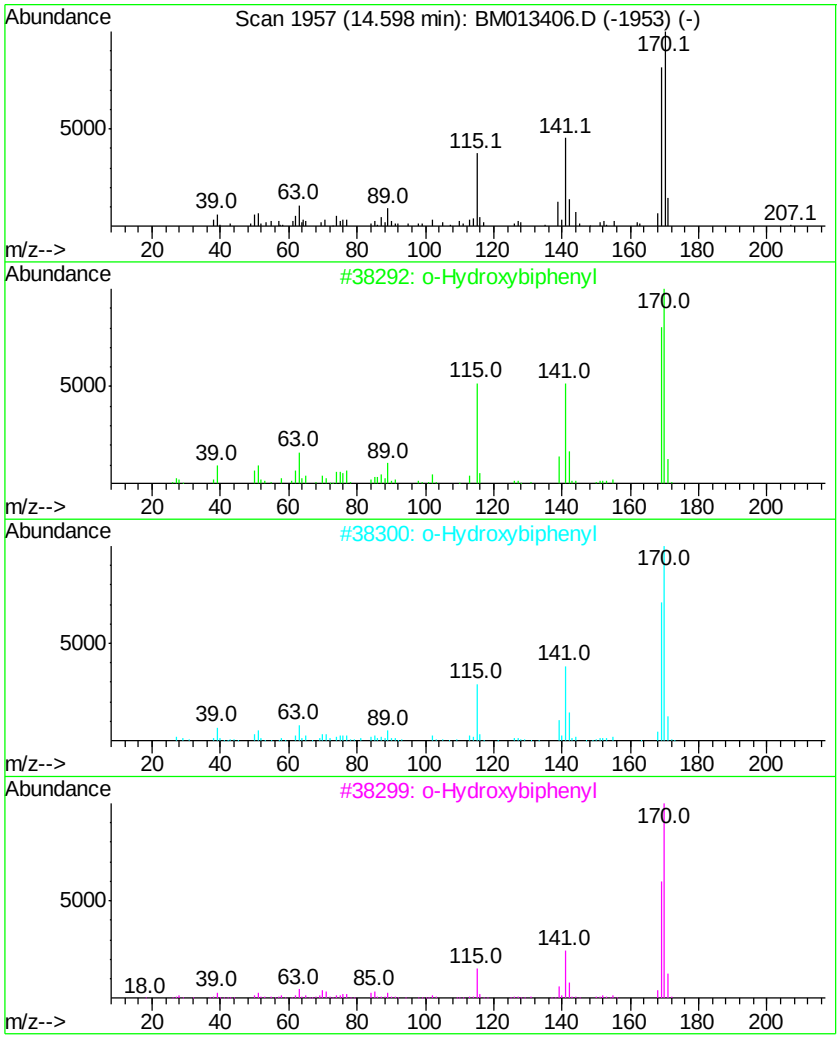
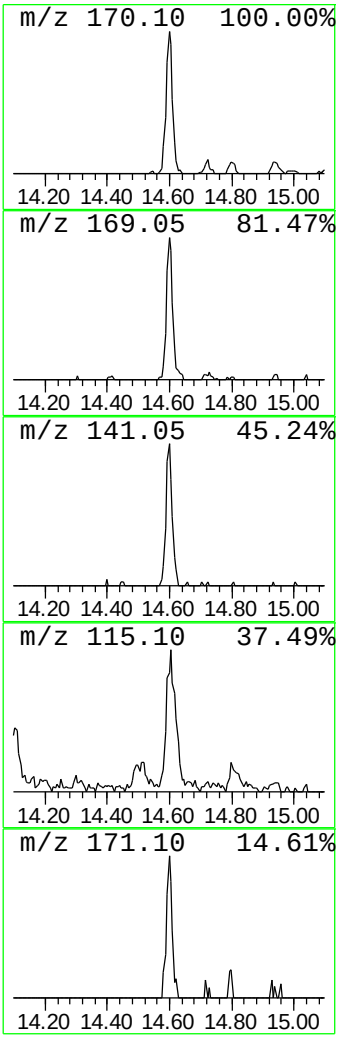
Peak Number 9 o-Hydroxybiphenyl Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.60	2.58 ng/ul	280953	Acenaphthene-d10	14.32

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	o-Hydroxybiphenyl	170	C12H10O	000090-43-7	97
2	o-Hydroxybiphenyl	170	C12H10O	000090-43-7	97
3	o-Hydroxybiphenyl	170	C12H10O	000090-43-7	94
4	o-Hydroxybiphenyl	170	C12H10O	000090-43-7	94
5	o-Hydroxybiphenyl	170	C12H10O	000090-43-7	94





Library Search Compound Report

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Acq On : 14 Dec 2017 17:34
Operator : SJ/JU
Sample : I6900-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : F9C00

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM113017.M
Quant Title : SVOA CALIBRATION

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

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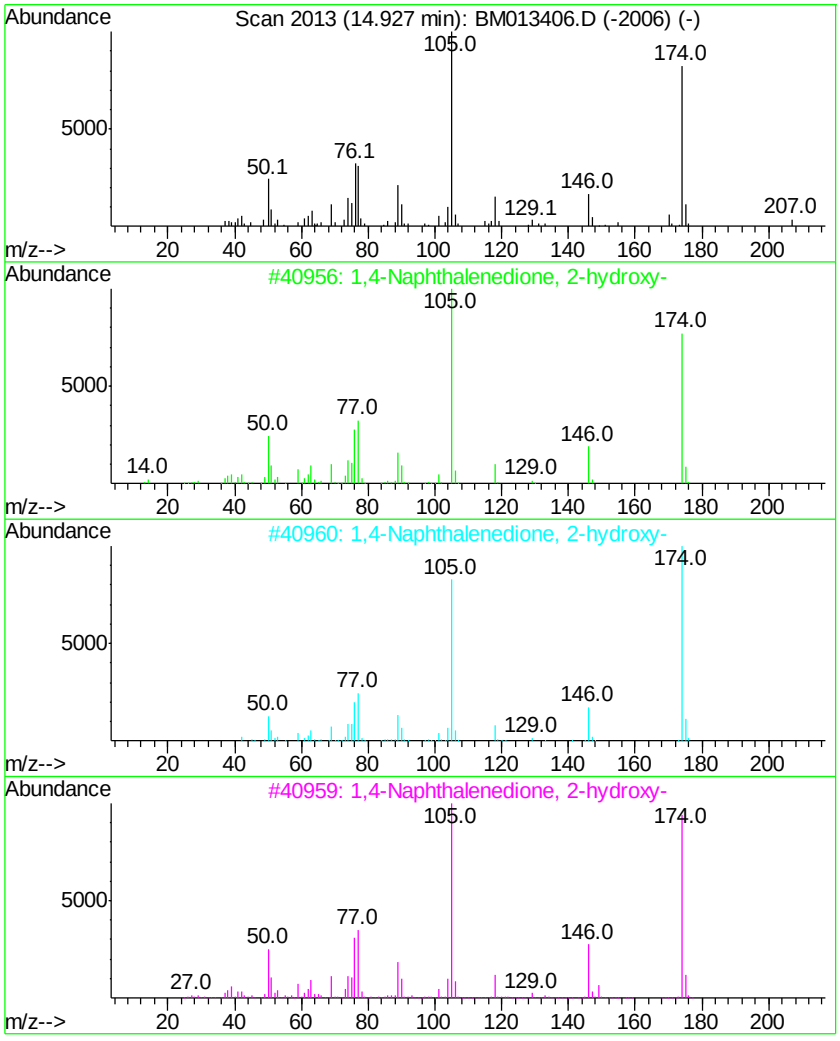
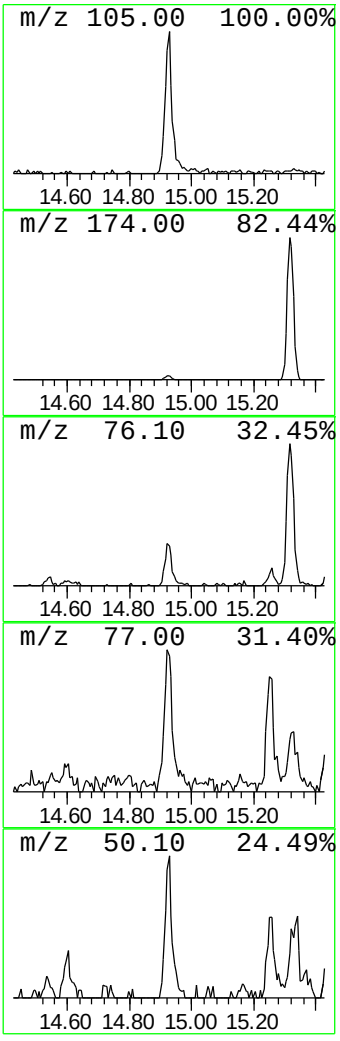
Peak Number 10 1,4-Naphthalenedione, 2-hyd... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.93	2.46 ng/ul	267675	Acenaphthene-d10	14.32

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,4-Naphthalenedione, 2-hydroxy-	174	C10H6O3	000083-72-7	97
2	1,4-Naphthalenedione, 2-hydroxy-	174	C10H6O3	000083-72-7	94
3	1,4-Naphthalenedione, 2-hydroxy-	174	C10H6O3	000083-72-7	93
4	Naphth[2,3-b]oxirene-2,7-dione, ...	174	C10H6O3	015448-58-5	90
5	1,4-Naphthalenedione, 2-hydroxy-	174	C10H6O3	000083-72-7	87





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Operator : SJ/JU
Sample : I6900-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : F9C00

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM113017.M
Quant Title : SVOA CALIBRATION

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

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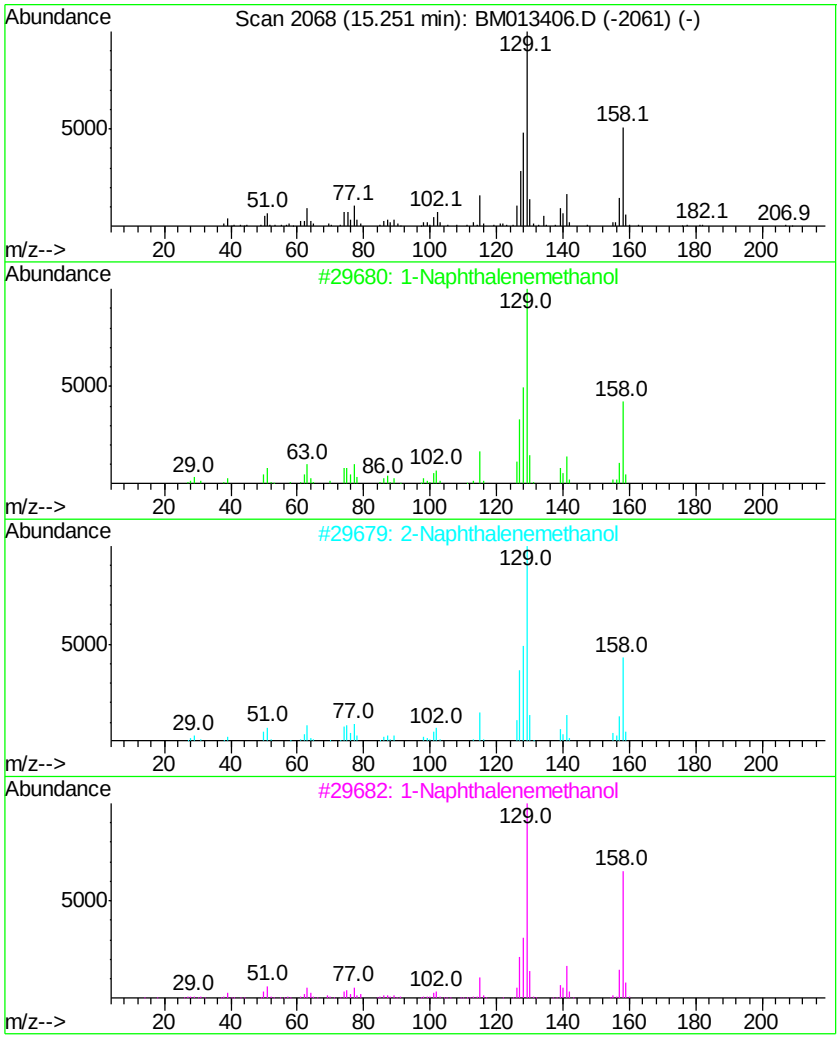
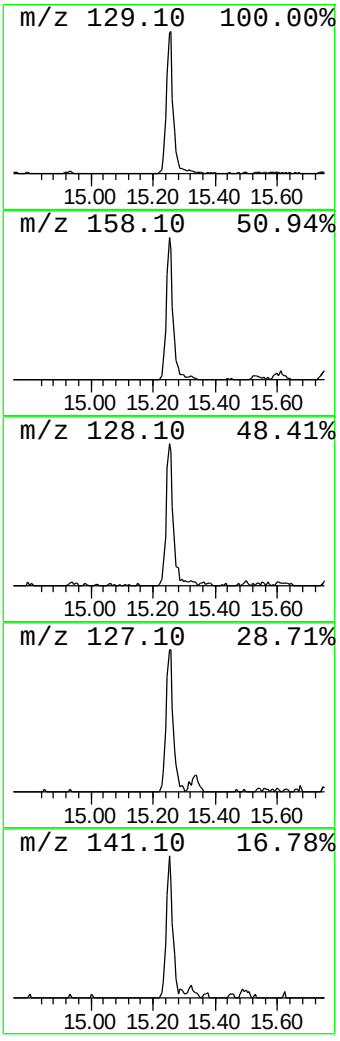
Peak Number 11 1-Naphthalenemethanol Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.25	4.80 ng/ul	522694	Acenaphthene-d10	14.32

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Naphthalenemethanol	158	C11H10O	004780-79-4	97
2	2-Naphthalenemethanol	158	C11H10O	001592-38-7	93
3	1-Naphthalenemethanol	158	C11H10O	004780-79-4	90
4	1,4-Methanonaphthalen-9-ol, 1,4-...	158	C11H10O	004796-33-2	87
5	1-Naphthalenemethanol	158	C11H10O	004780-79-4	87





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Acq On : 14 Dec 2017 17:34
Operator : SJ/JU
Sample : I6900-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : F9C00

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM113017.M
Quant Title : SVOA CALIBRATION

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

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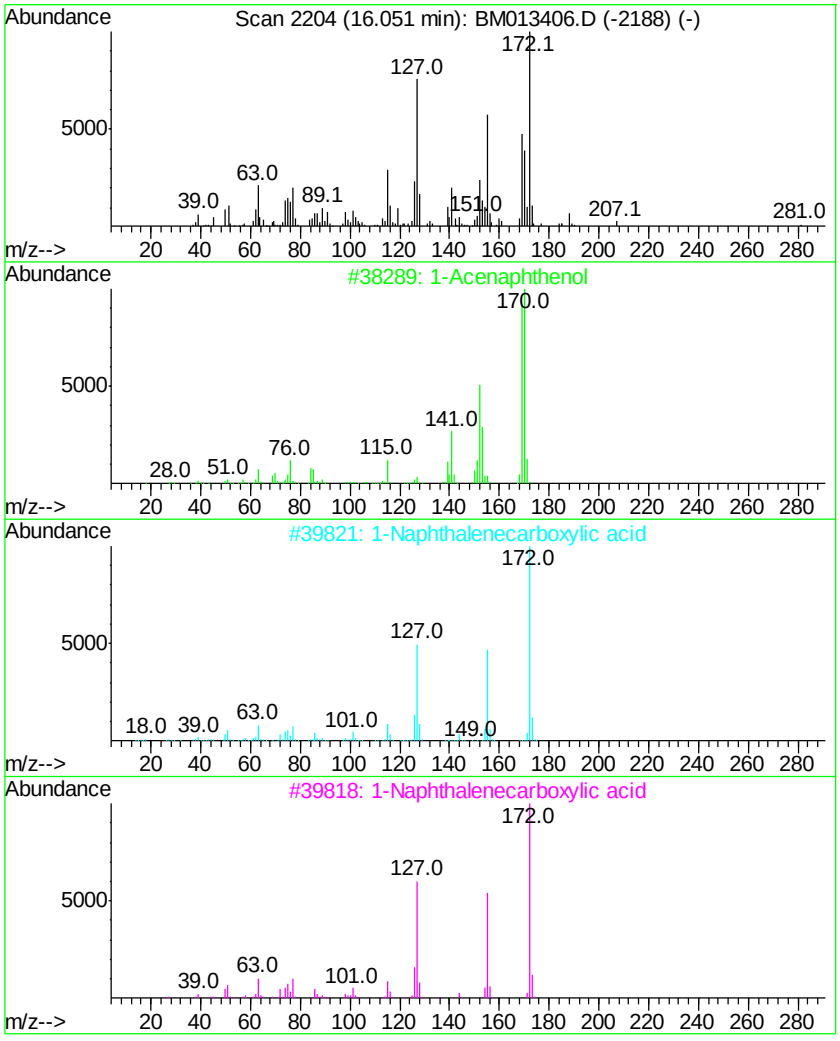
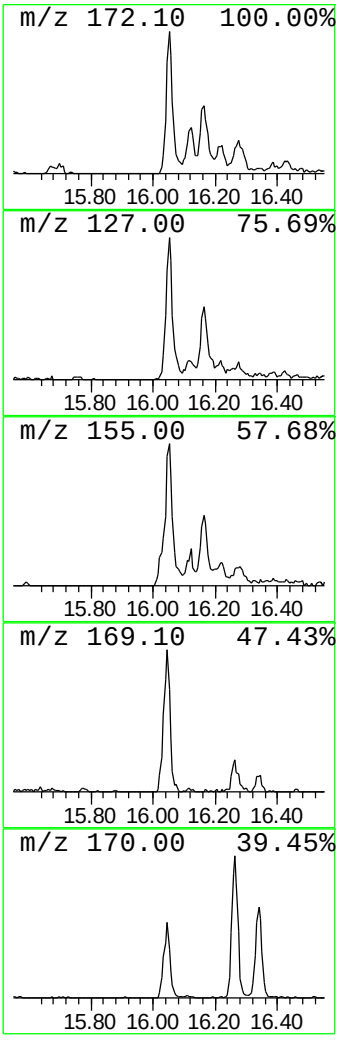
Peak Number 12 1-Acenaphthenol Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.05	5.19 ng/ul	791957	Phenanthrene-d10	17.07

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Acenaphthenol	170	C12H10O	006306-07-6	50
2	1-Naphthalenecarboxylic acid	172	C11H8O2	000086-55-5	49
3	1-Naphthalenecarboxylic acid	172	C11H8O2	000086-55-5	49
4	o-Hydroxybiphenyl	170	C12H10O	000090-43-7	47
5	2-Naphthyl methyl ketone	170	C12H10O	000093-08-3	42





Library Search Compound Report

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Acq On : 14 Dec 2017 17:34
Operator : SJ/JU
Sample : I6900-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : F9C00

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM113017.M
Quant Title : SVOA CALIBRATION

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

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Peak Number 13 p-Hydroxybiphenyl Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.26	2.13 ng/ul	324303	Phenanthrene-d10	17.07

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	p-Hydroxybiphenyl	170	C12H10O	000092-69-3	91
2	p-Hydroxybiphenyl	170	C12H10O	000092-69-3	91
3	3-Hydroxybiphenyl	170	C12H10O	000580-51-8	87
4	p-Hydroxybiphenyl	170	C12H10O	000092-69-3	87
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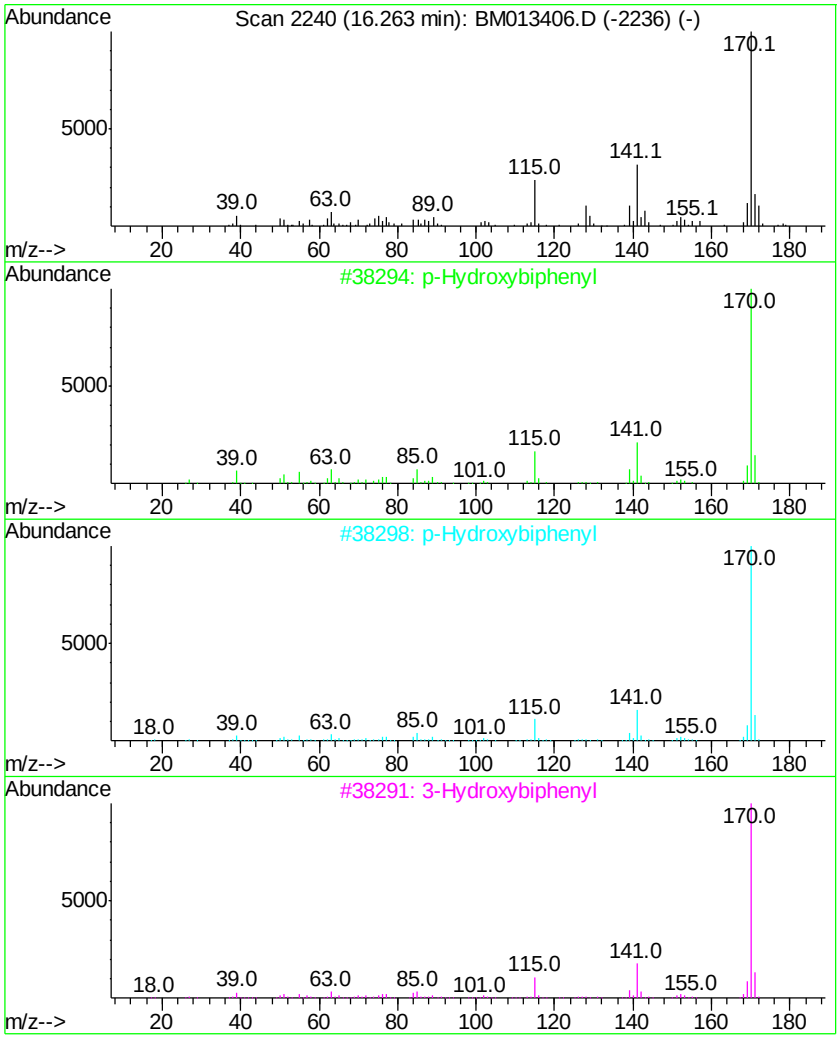

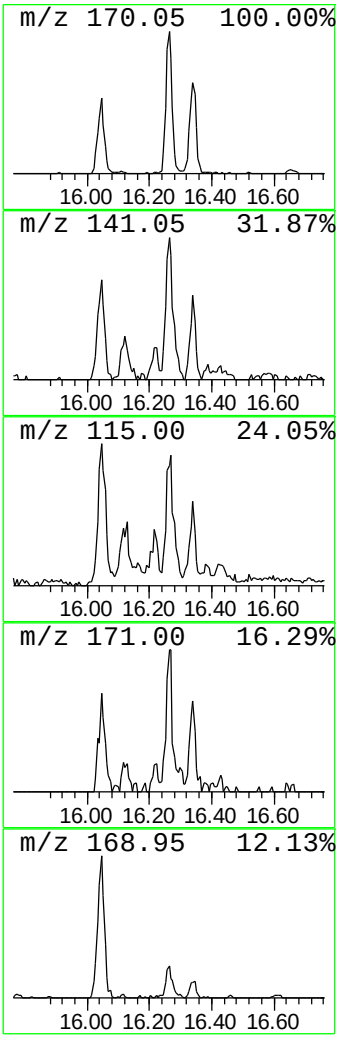

Tentatively Identified Compound (LSC) summary

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Acq On : 14 Dec 2017 17:34
Operator : SJ/JU
Sample : I6900-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : F9C00

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM113017.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
Cyclohexanone	5.75	4.2	ng/ul	218948	1	7.68	1035490	20.0
Phenol, 2,6-dimet...	9.09	3.0	ng/ul	220929	2	10.46	1469940	20.0
Phenol, 3,5-dimet...	9.95	18.6	ng/ul	1369030	2	10.46	1469940	20.0
3-Methylcyclohexa...	9.99	5.7	ng/ul	418166	2	10.46	1469940	20.0
Benzoic acid, 3-m...	11.26	4.9	ng/ul	360070	2	10.46	1469940	20.0
Phenol, 2-ethyl-6...	11.29	4.0	ng/ul	291949	2	10.46	1469940	20.0
Phenol, 4-(2-prop...	12.23	3.6	ng/ul	265844	2	10.46	1469940	20.0
Cyclohexene, 3-me...	13.85	7.7	ng/ul	843259	3	14.32	2178750	20.0
o-Hydroxybiphenyl	14.60	2.6	ng/ul	280953	3	14.32	2178750	20.0
1,4-Naphthalenedi...	14.93	2.5	ng/ul	267675	3	14.32	2178750	20.0
1-Naphthalenemeth...	15.25	4.8	ng/ul	522694	3	14.32	2178750	20.0
1-Acenaphthenol	16.05	5.2	ng/ul	791957	4	17.07	3050270	20.0
p-Hydroxybiphenyl	16.26	2.1	ng/ul	324303	4	17.07	3050270	20.0