LSC Area Percent Report

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM061217\
Data File : BM010547.D
Acq On : 12 Jun 2017 22:21
Operator : SJ/MA
Sample : I3558-14DL 2X
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : F5C87DL

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-BM052717.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	4.010	147	157	165	rVB2	92340	205320	2.13%	0.377%
2	7.075	671	678	688	rBV2	250517	554565	5.76%	1.018%
3	7.228	698	704	711	rBV	165445	266120	2.76%	0.488%
4	7.428	730	738	747	rBV2	303657	510446	5.30%	0.937%
5	7.886	808	816	825	rBV	521396	815728	8.47%	1.497%
6	8.345	888	894	899	rBV3	74625	111717	1.16%	0.205%
7	8.610	932	939	945	rBV2	321001	548490	5.70%	1.007%
8	8.675	945	950	958	rVB3	135632	250928	2.61%	0.461%
9	9.075	1010	1018	1025	rBV2	232829	404257	4.20%	0.742%
10	9.786	1132	1139	1146	rBV2	223890	382413	3.97%	0.702%
11	9.869	1146	1153	1156	rBV2	114786	192954	2.00%	0.354%
12	9.986	1162	1173	1179	rBV	198047	489662	5.08%	0.899%
13	10.175	1201	1205	1213	rVB2	98311	160464	1.67%	0.295%
14	10.316	1222	1229	1238	rBV2	419883	741193	7.70%	1.360%
15	10.686	1285	1292	1296	rVV	598683	1041370	10.81%	1.911%
16	10.739	1296	1301	1308	rVB	1392068	2308031	23.97%	4.236%
17	10.857	1311	1321	1329	rBV	362123	695478	7.22%	1.276%
18	11.251	1379	1388	1395	rBV4	82949	205900	2.14%	0.378%
19	11.522	1429	1434	1438	rVB2	79786	130572	1.36%	0.240%
20	12.339	1566	1573	1579	rBV	299321	509234	5.29%	0.935%
21	12.421	1581	1587	1594	rBV2	398065	624171	6.48%	1.146%
22	12.563	1605	1611	1617	rVB2	162493	276207	2.87%	0.507%
23	12.698	1629	1634	1642	rBV4	133175	227101	2.36%	0.417%
24	13.357	1740	1746	1752	rVB2	78385	130408	1.35%	0.239%
25	13.680	1795	1801	1809	rVB7	50970	143765	1.49%	0.264%
26	13.939	1840	1845	1849	rVV	752313	1031711	10.71%	1.894%
27	13.986	1849	1853	1862	rVB	332705	466094	4.84%	0.855%
28	14.221	1887	1893	1903	rVB	732436	1127378	11.71%	2.069%
29	14.374	1913	1919	1929	rBV	165207	293751	3.05%	0.539%
30	14.521	1939	1944	1950	rVV2	958309	1491018	15.48%	2.736%
31	14.586	1950	1955	1962	rVB2	302918	444141	4.61%	0.815%
32	14.774	1980	1987	1995	rBV2	209560	461804	4.80%	0.848%
33	14.839	1995	1998	2006	rVV3	68120	112448	1.17%	0.206%
34	14.921	2006	2012	2017	rVV	293215	455592	4.73%	0.836%

LSC Area Percent Report

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Data File : BM010547.D
Acq On : 12 Jun 2017 22:21
Operator : SJ/MA
Sample : I3558-14DL 2X
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : F5C87DL

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-BM052717.M
Title : SVOA CALIBRATION

35	15.280	2068	2073	2081	rBV7	46923	110473	1.15%	0.203%
36	15.515	2102	2113	2118	rBV	1037840	1506866	15.65%	2.766%
37	15.568	2118	2122	2127	rVB	370897	542334	5.63%	0.995%
38	15.645	2130	2135	2141	rBV	245712	370657	3.85%	0.680%
39	16.004	2192	2196	2205	rVB	72525	163841	1.70%	0.301%
40	16.551	2264	2289	2292	rBV3	2258912	9630250	100.00%	17.675%
41	16.845	2334	2339	2345	rVB4	79393	149756	1.56%	0.275%
42	17.092	2372	2381	2385	rBV	101589	184341	1.91%	0.338%
43	17.274	2406	2412	2415	rBV	1525387	2093623	21.74%	3.842%
44	17.315	2415	2419	2424	rVV	1454453	1939339	20.14%	3.559%
45	17.368	2424	2428	2432	rVV	1251925	1911946	19.85%	3.509%
46	17.404	2432	2434	2442	rVB	426674	627443	6.52%	1.152%
47	17.692	2478	2483	2493	rVB	323276	505657	5.25%	0.928%
48	17.792	2493	2500	2506	rBV6	54494	124567	1.29%	0.229%
49	18.127	2551	2557	2563	rBV3	172825	372364	3.87%	0.683%
50	18.186	2563	2567	2573	rVB	131169	198486	2.06%	0.364%
51	18.339	2587	2593	2597	rBV2	164243	256351	2.66%	0.470%
52	18.639	2640	2644	2649	rVB	94638	134227	1.39%	0.246%
53	19.321	2755	2760	2766	rVB	915324	1221842	12.69%	2.242%
54	19.392	2766	2772	2775	rBV7	70340	108787	1.13%	0.200%
55	19.527	2790	2795	2800	rBV	1220994	1384256	14.37%	2.541%
56	19.662	2812	2818	2821	rBV	1705381	2610641	27.11%	4.791%
57	20.174	2902	2905	2912	rVB	128047	195658	2.03%	0.359%
58	20.274	2918	2922	2927	rBV	134788	196366	2.04%	0.360%
59	20.568	2968	2972	2976	rBV2	769758	821239	8.53%	1.507%
60	21.092	3058	3061	3065	rBV6	115994	188441	1.96%	0.346%
61	21.439	3114	3120	3124	rBV	2656303	3489997	36.24%	6.405%
62	21.474	3124	3126	3130	rVB	180219	207026	2.15%	0.380%
63	23.621	3485	3491	3497	rBV2	1310690	2426351	25.20%	4.453%
64	23.774	3511	3517	3524	rVB	1656145	3003060	31.18%	5.512%

Sum of corrected areas: 54486616

LSC Report - Integrated Chromatogram

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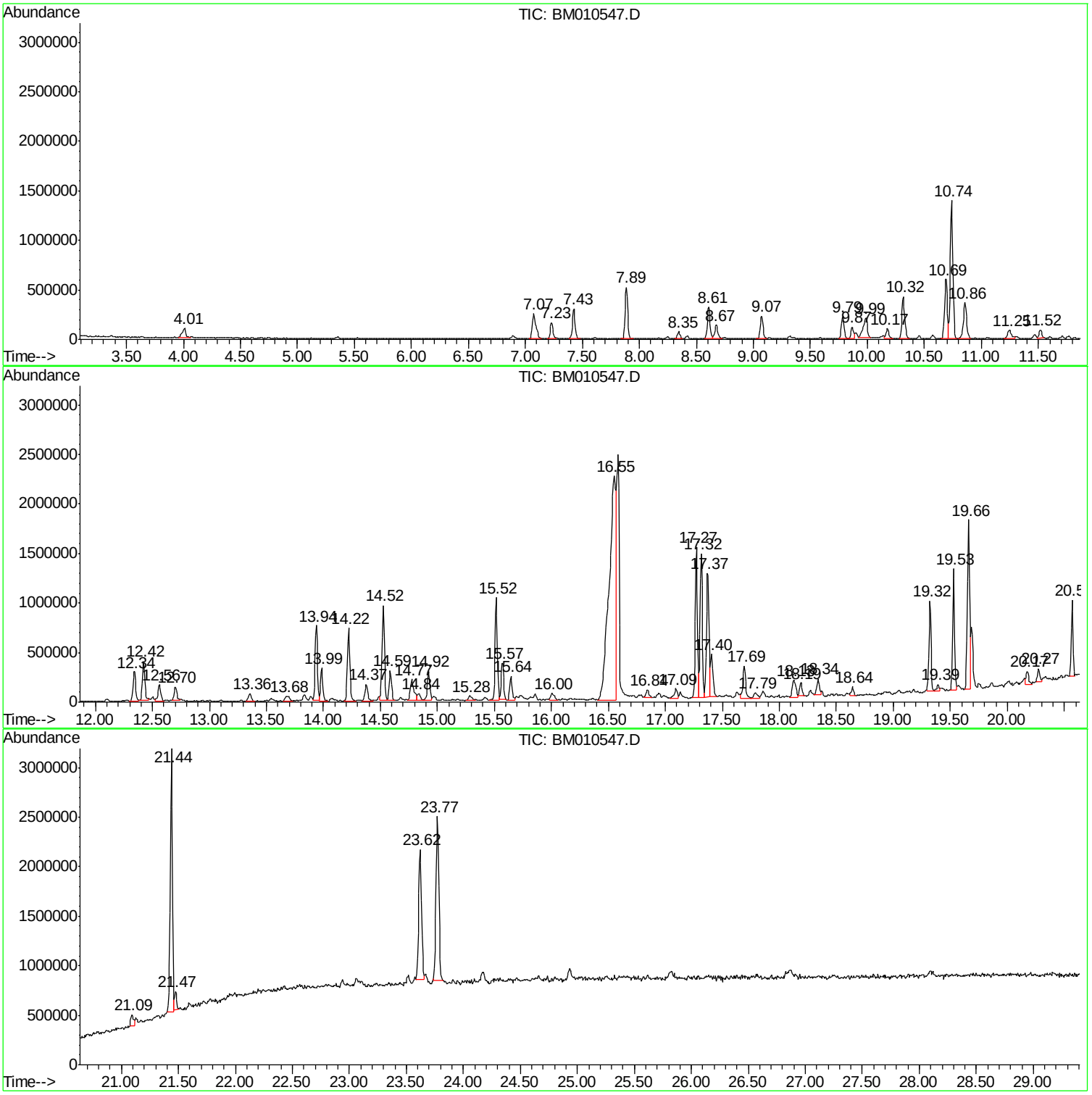
Data File : BM010547.D
Acq On : 12 Jun 2017 22:21
Operator : SJ/MA
Sample : I3558-14DL 2X
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : F5C87DL

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM052717.M
Quant Title : SVOA CALIBRATION

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



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ALS Vial : 21 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : F5C87DL

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM052717.M
Quant Title : SVOA CALIBRATION

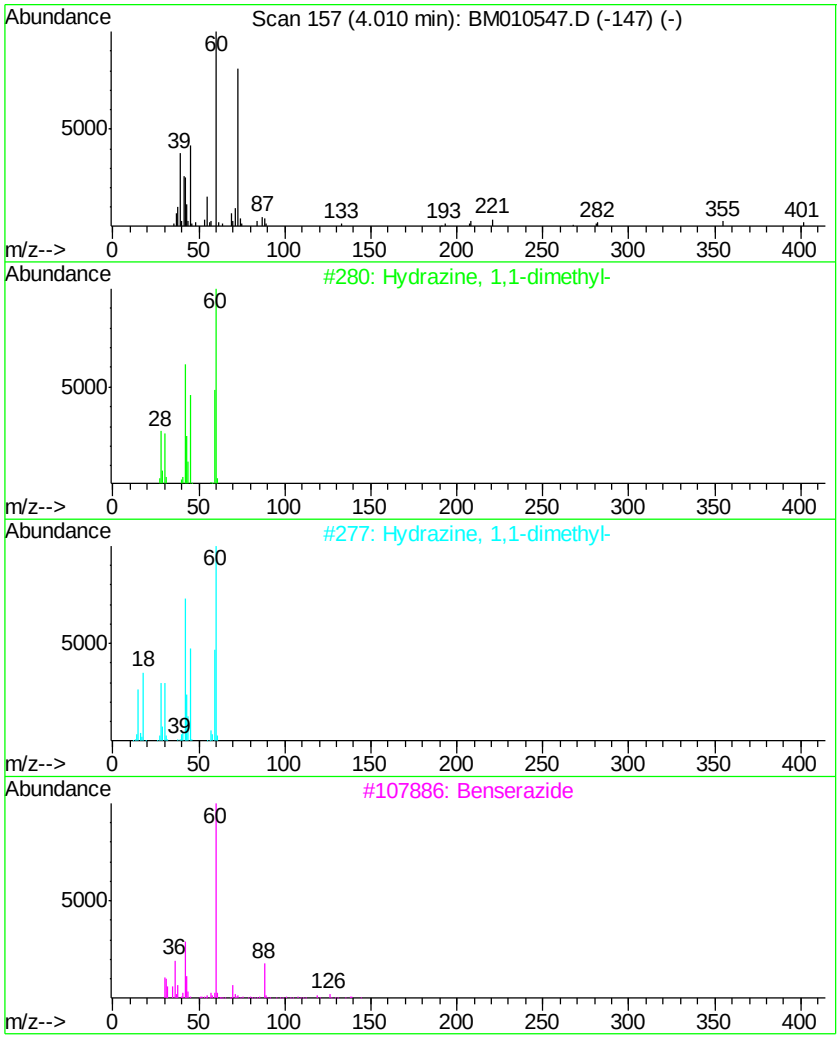
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

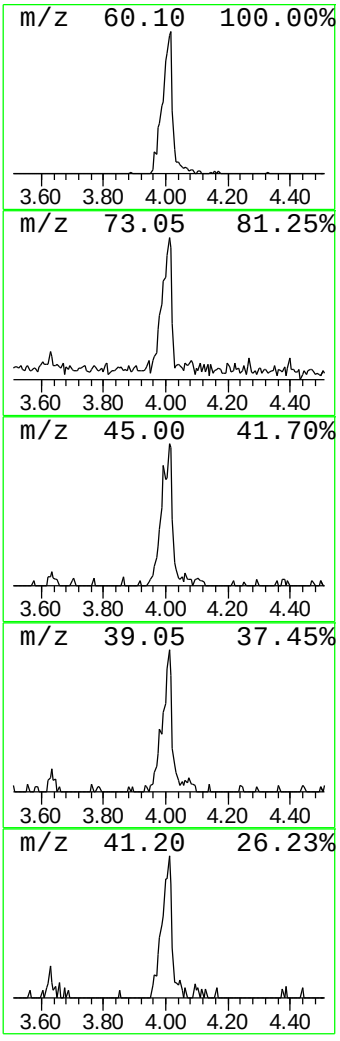
****

Peak Number 1 unknown-01 Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.01	5.03 ng/ul	205320	1,4-Dichlorobenzene-d4	7.89

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hydrazine, 1,1-dimethyl-	60	C2H8N2	000057-14-7	40
2	Hydrazine, 1,1-dimethyl-	60	C2H8N2	000057-14-7	33
3	Benserazide	257	C10H15N3O5	000322-35-0	25
4	Hydrazine, 1,1-diethyl-	88	C4H12N2	000616-40-0	25
5	1-Butanol, 3-(1-ethoxyethoxy)-2-...	176	C9H20O3	059410-44-5	12





Data Path : Z:\HPCHEM1\BNA_M\DATA\BM061217\
Data File : BM010547.D
Acq On : 12 Jun 2017 22:21
Operator : SJ/MA
Sample : I3558-14DL 2X
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : F5C87DL

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM052717.M
Quant Title : SVOA CALIBRATION

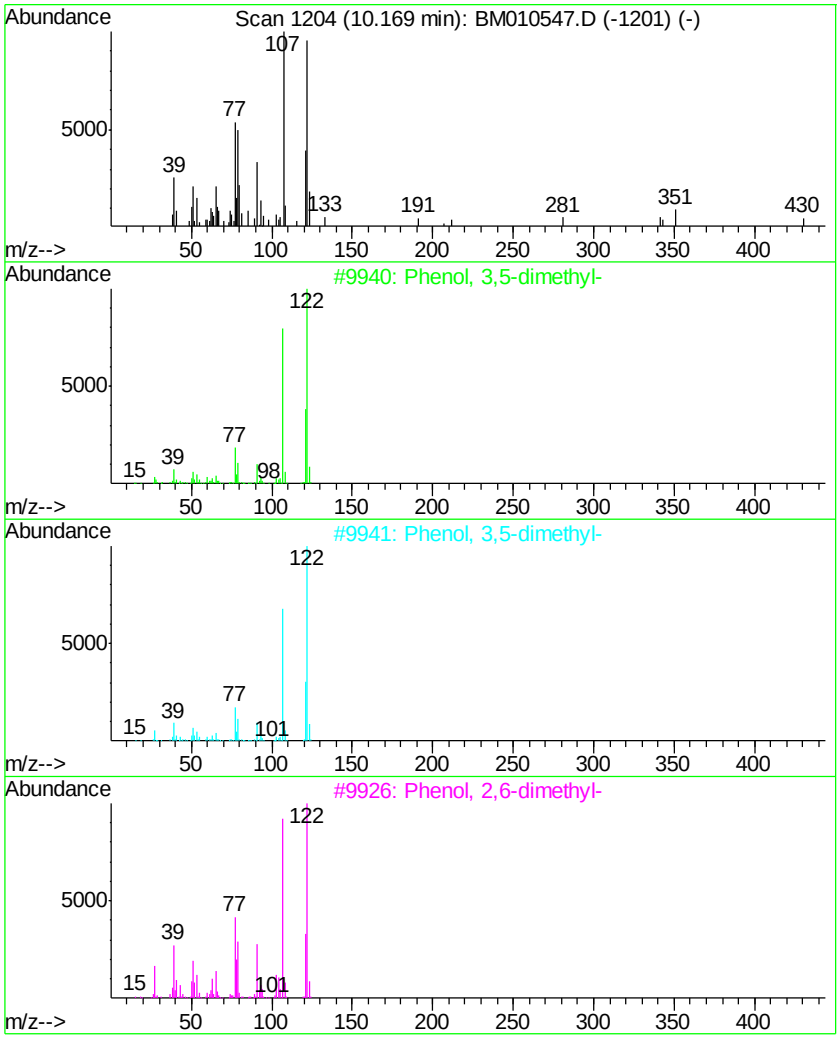
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

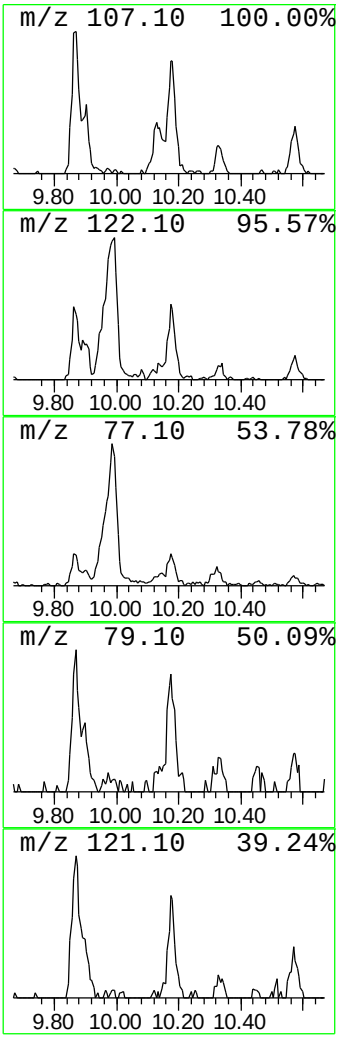
****

Peak Number 3 Phenol, 3,5-dimethyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.17	3.08 ng/ul	160464	Naphthalene-d8	10.69

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, 3,5-dimethyl-	122	C8H10O	000108-68-9	91
2	Phenol, 3,5-dimethyl-	122	C8H10O	000108-68-9	90
3	Phenol, 2,6-dimethyl-	122	C8H10O	000576-26-1	87
4	Phenol, 2,6-dimethyl-	122	C8H10O	000576-26-1	86
5	Phenol, 3,5-dimethyl-	122	C8H10O	000108-68-9	86





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Sample : I3558-14DL 2X
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : F5C87DL

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM052717.M
Quant Title : SVOA CALIBRATION

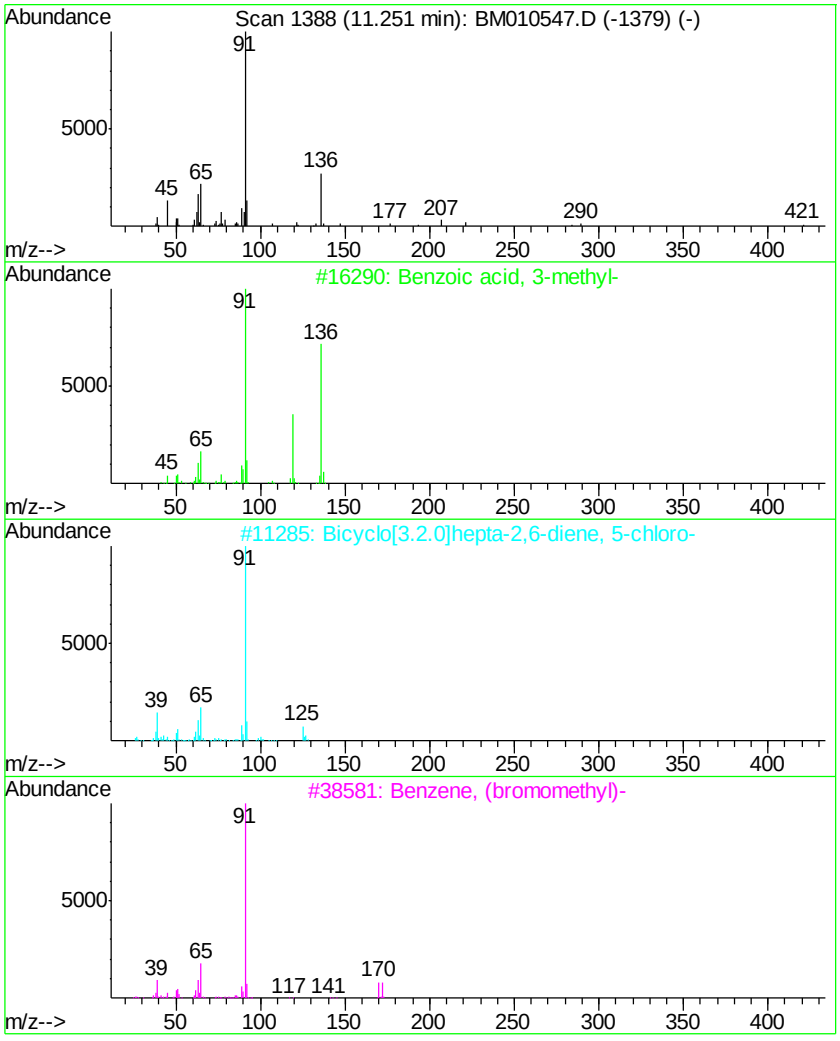
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

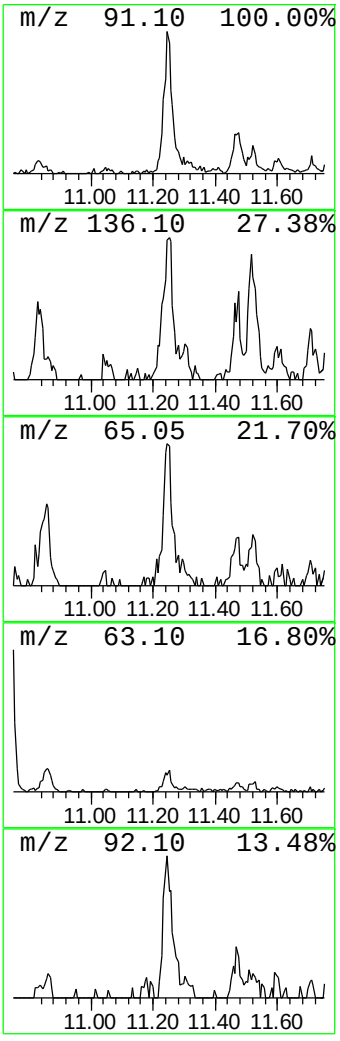
****

Peak Number 4 Benzoic acid, 3-methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.25	3.95 ng/ul	205900	Naphthalene-d8	10.69

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzoic acid, 3-methyl-	136	C8H8O2	000099-04-7	64
2	Bicyclo[3.2.0]hepta-2,6-diene, 5...	126	C7H7Cl	023610-81-3	53
3	Benzene, (bromomethyl)-	170	C7H7Br	000100-39-0	53
4	3,5,6-Trichloro-2-pyridyl .alpha...	351	C12H8Cl3NO3S	058236-69-4	50
5	2,3,5-Trioxabicyclo[2.1.0]pentan...	254	C16H14O3	056247-48-4	50





Data Path : Z:\HPCHEM1\BNA_M\DATA\BM061217\
Data File : BM010547.D
Acq On : 12 Jun 2017 22:21
Operator : SJ/MA
Sample : I3558-14DL 2X
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : F5C87DL

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM052717.M
Quant Title : SVOA CALIBRATION

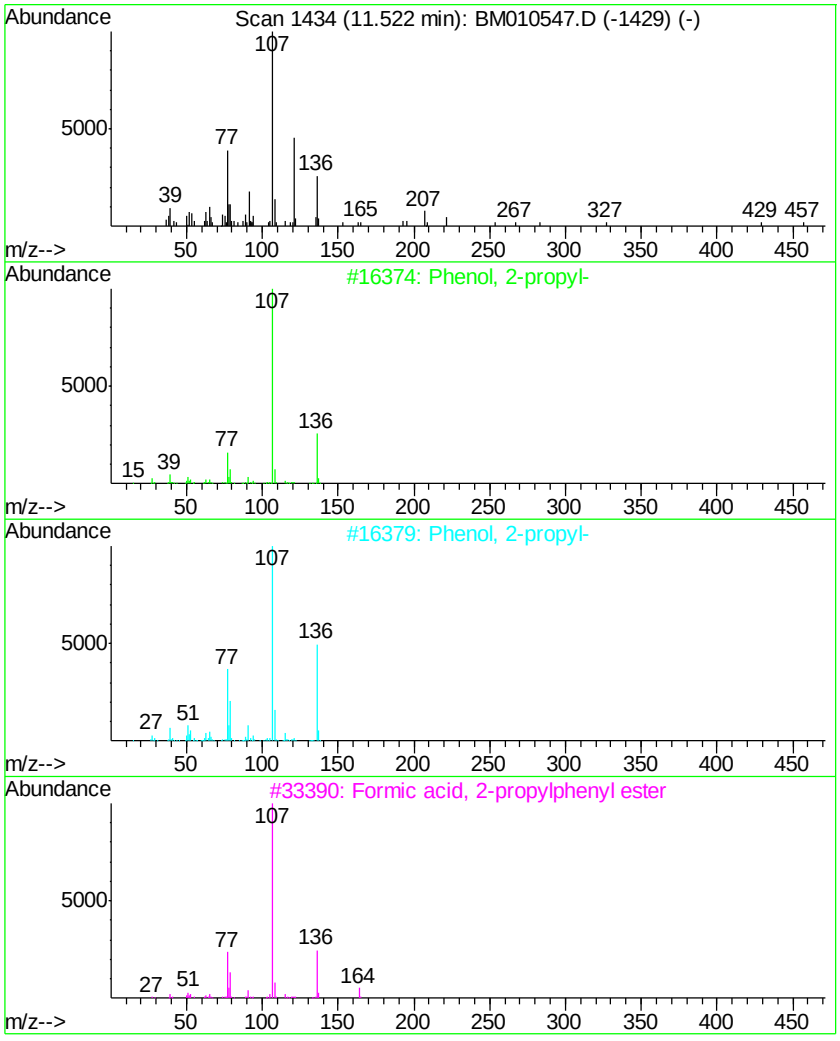
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

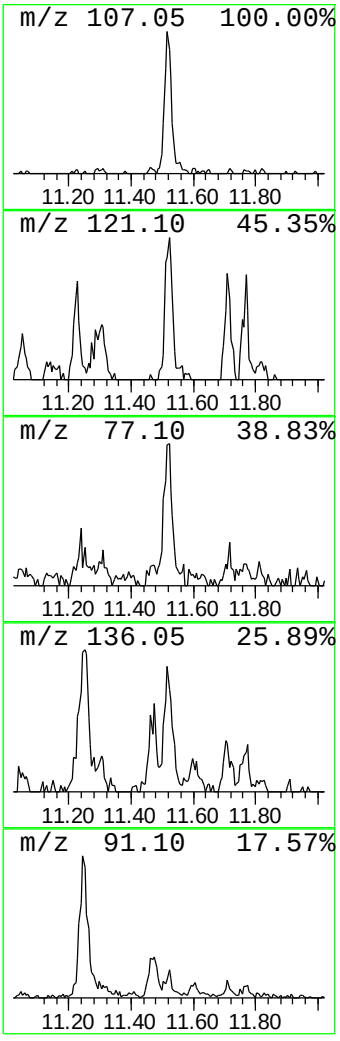
****

Peak Number 5 Phenol, 2-propyl- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.52	2.51 ng/ul	130572	Naphthalene-d8	10.69

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, 2-propyl-	136	C9H12O	000644-35-9	72
2	Phenol, 2-propyl-	136	C9H12O	000644-35-9	72
3	Formic acid, 2-propylphenyl ester	164	C10H12O2	1000368-96-6	59
4	d-Tyrosine	181	C9H11NO3	000556-02-5	59
5	Phenol, 4-ethyl-	122	C8H10O	000123-07-9	59





Library Search Compound Report

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Acq On : 12 Jun 2017 22:21
Operator : SJ/MA
Sample : I3558-14DL 2X
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : F5C87DL

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM052717.M
Quant Title : SVOA CALIBRATION

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

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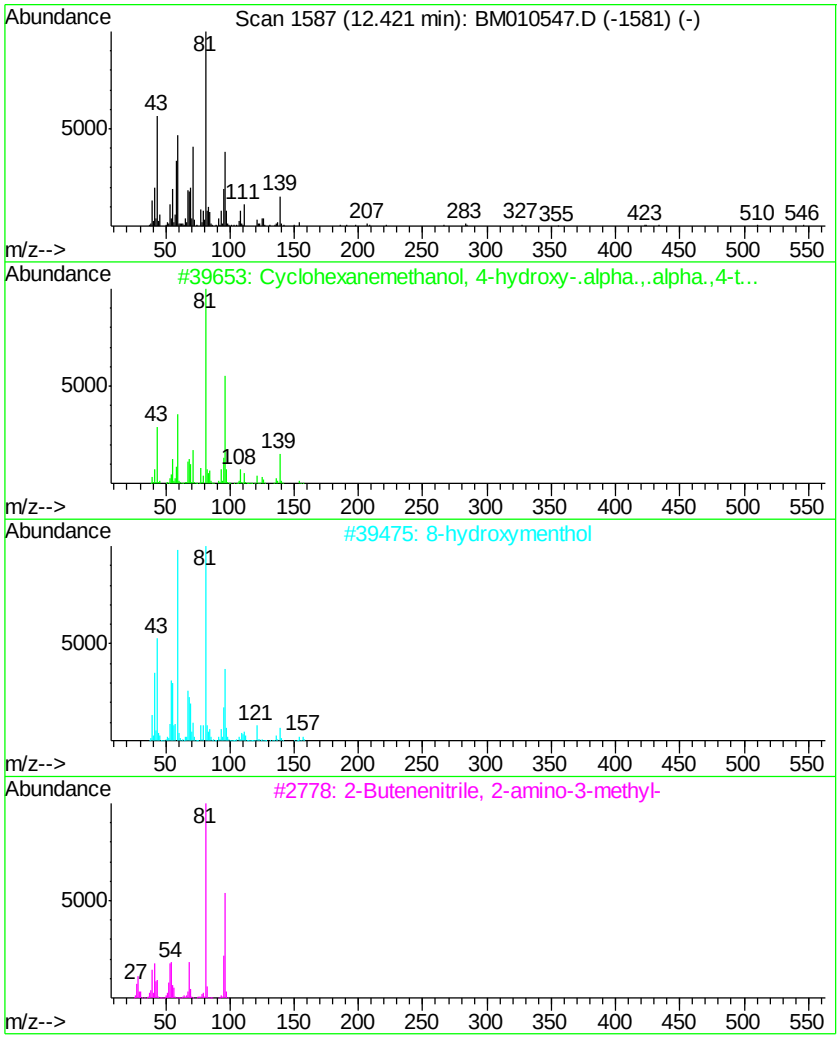
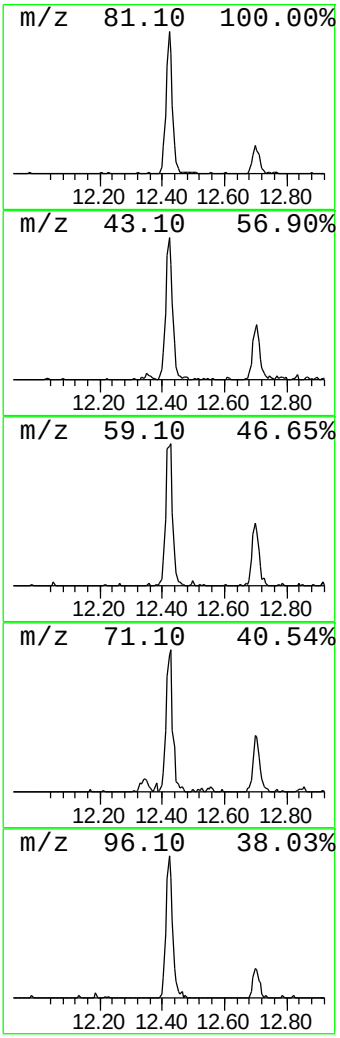
Peak Number 6 Cyclohexanemethanol, 4-hydr... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.42	11.99 ng/ul	624171	Naphthalene-d8	10.69

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclohexanemethanol, 4-hydroxy-....	172	C10H20O2	000080-53-5	53
2	8-hydroxymenthol	172	C10H20O2	1000374-16-2	53
3	2-Butenenitrile, 2-amino-3-methyl-	96	C5H8N2	1000196-28-0	38
4	3,5-Dimethylcyclopentene	96	C7H12	007459-71-4	35
5	1,4-Hexadiene, 4-methyl-	96	C7H12	001116-90-1	27





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Acq On : 12 Jun 2017 22:21
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Sample : I3558-14DL 2X
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : F5C87DL

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM052717.M
Quant Title : SVOA CALIBRATION

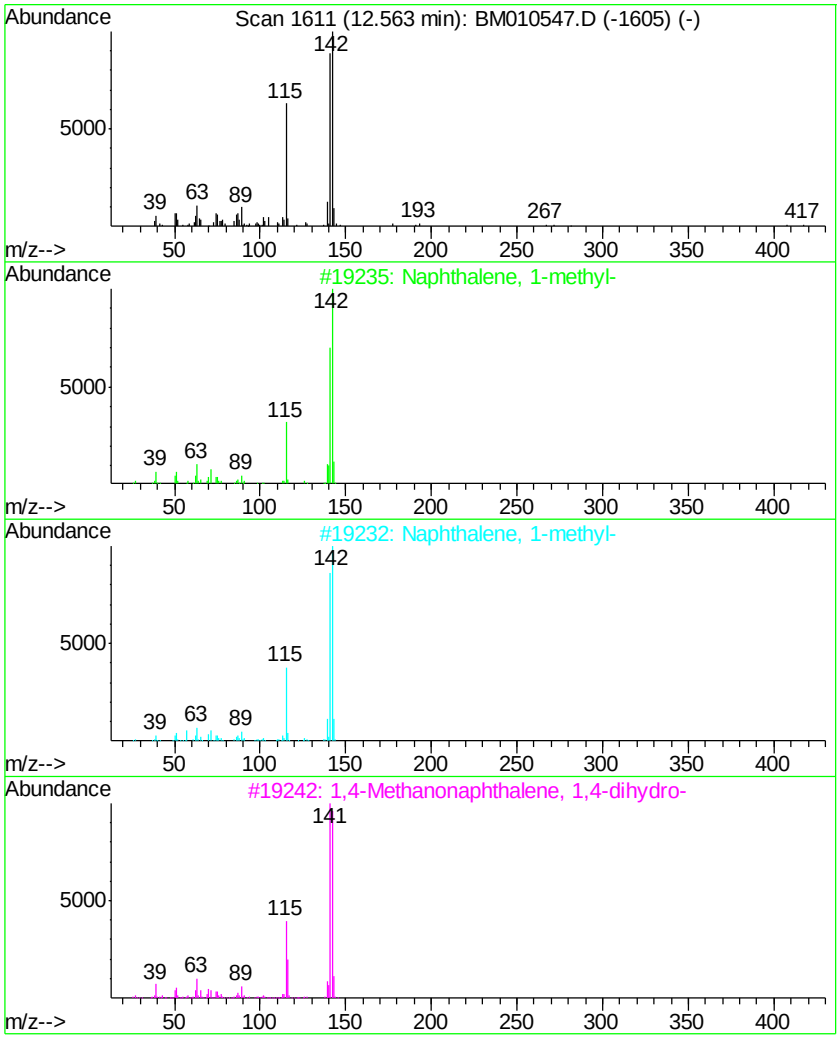
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

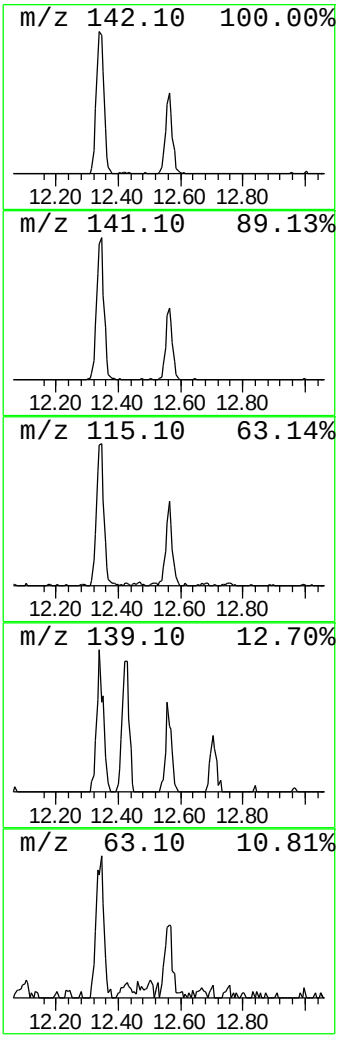
****

Peak Number 7 Naphthalene, 1-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.56	5.30 ng/ul	276207	Naphthalene-d8	10.69

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 1-methyl-	142	C11H10	000090-12-0	94
2	Naphthalene, 1-methyl-	142	C11H10	000090-12-0	94
3	1,4-Methanonaphthalene, 1,4-dihy...	142	C11H10	004453-90-1	91
4	Naphthalene, 2-methyl-	142	C11H10	000091-57-6	91
5	Naphthalene, 2-methyl-	142	C11H10	000091-57-6	91





Library Search Compound Report

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Data File : BM010547.D
Acq On : 12 Jun 2017 22:21
Operator : SJ/MA
Sample : I3558-14DL 2X
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : F5C87DL

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM052717.M
Quant Title : SVOA CALIBRATION

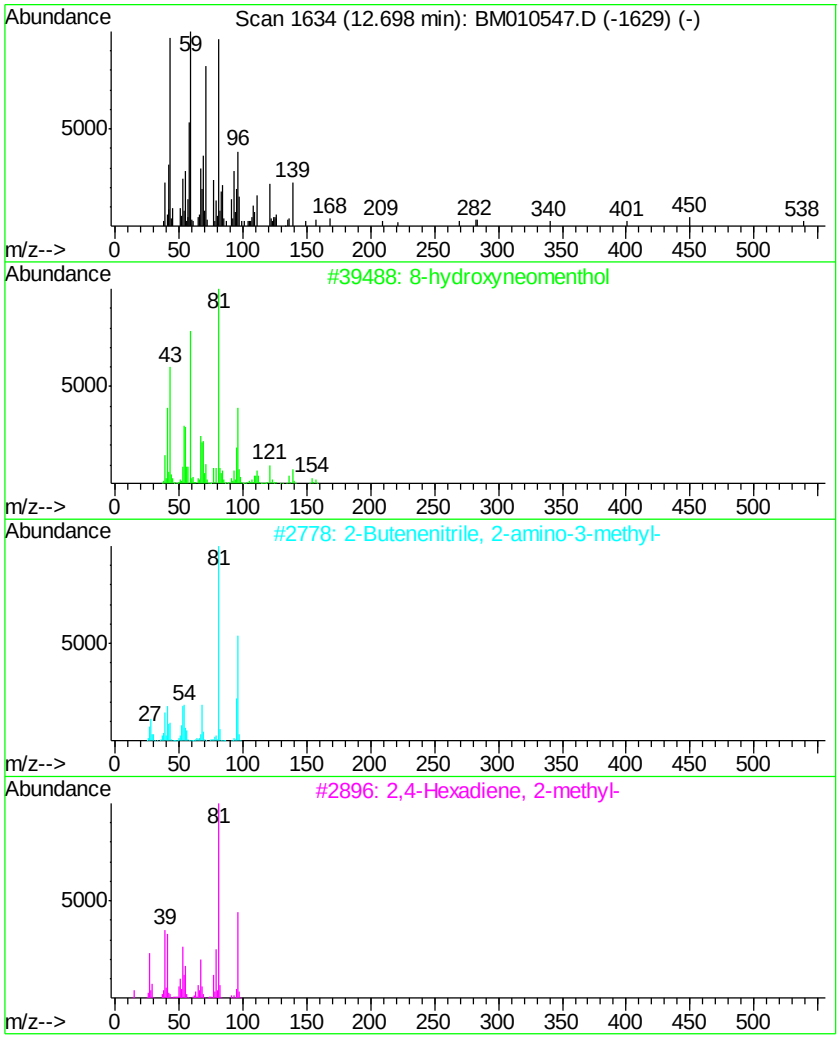
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

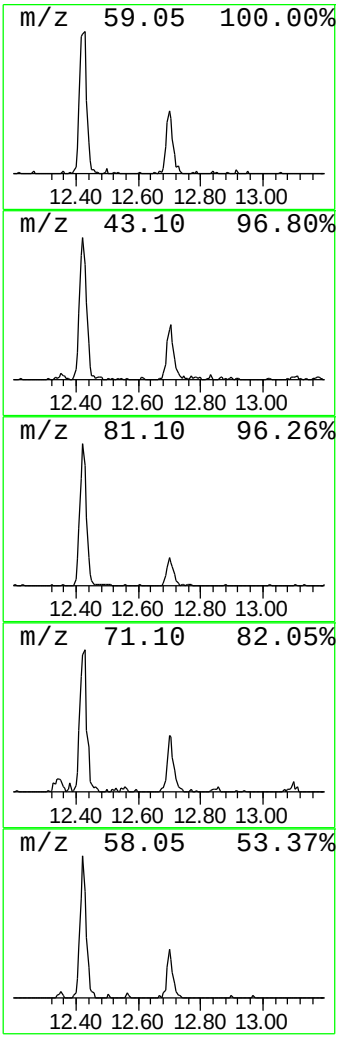
****

Peak Number 8 unknown-02 Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.70	3.05 ng/ul	227101	Acenaphthene-d10	14.52

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	8-hydroxyneomenthol	172	C10H20O2	1000374-16-1	47
2	2-Butenenitrile, 2-amino-3-methyl-	96	C5H8N2	1000196-28-0	25
3	2,4-Hexadiene, 2-methyl-	96	C7H12	028823-41-8	18
4	Cyclohexene, 1-methyl-	96	C7H12	000591-49-1	18
5	Carbonic acid, propargyl cyclohe...	196	C11H16O3	1000357-91-0	18





Library Search Compound Report

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM061217\
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Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : F5C87DL

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM052717.M
Quant Title : SVOA CALIBRATION

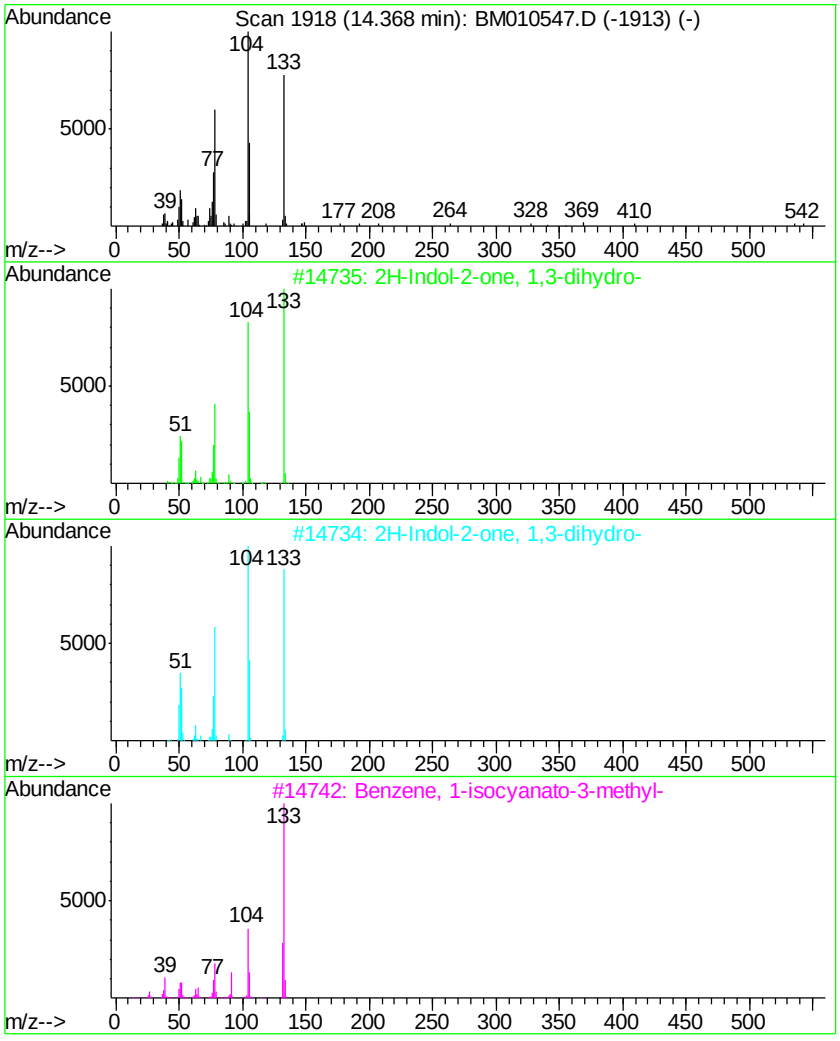
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

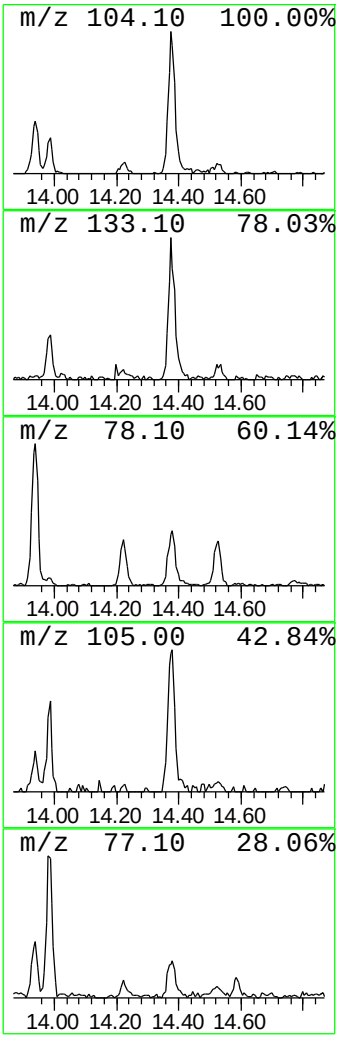
****

Peak Number 9 2H-Indol-2-one, 1,3-dihydro- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.37	3.94 ng/ul	293751	Acenaphthene-d10	14.52

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2H-Indol-2-one, 1,3-dihydro-	133	C8H7NO	000059-48-3	91
2	2H-Indol-2-one, 1,3-dihydro-	133	C8H7NO	000059-48-3	91
3	Benzene, 1-isocyanato-3-methyl-	133	C8H7NO	000621-29-4	91
4	1H-Benzotriazole, 4-methyl-	133	C7H7N3	029878-31-7	87
5	Benzene, 1-isocyanato-3-methyl-	133	C8H7NO	000621-29-4	86





Library Search Compound Report

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Data File : BM010547.D
Acq On : 12 Jun 2017 22:21
Operator : SJ/MA
Sample : I3558-14DL 2X
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : F5C87DL

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM052717.M
Quant Title : SVOA CALIBRATION

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

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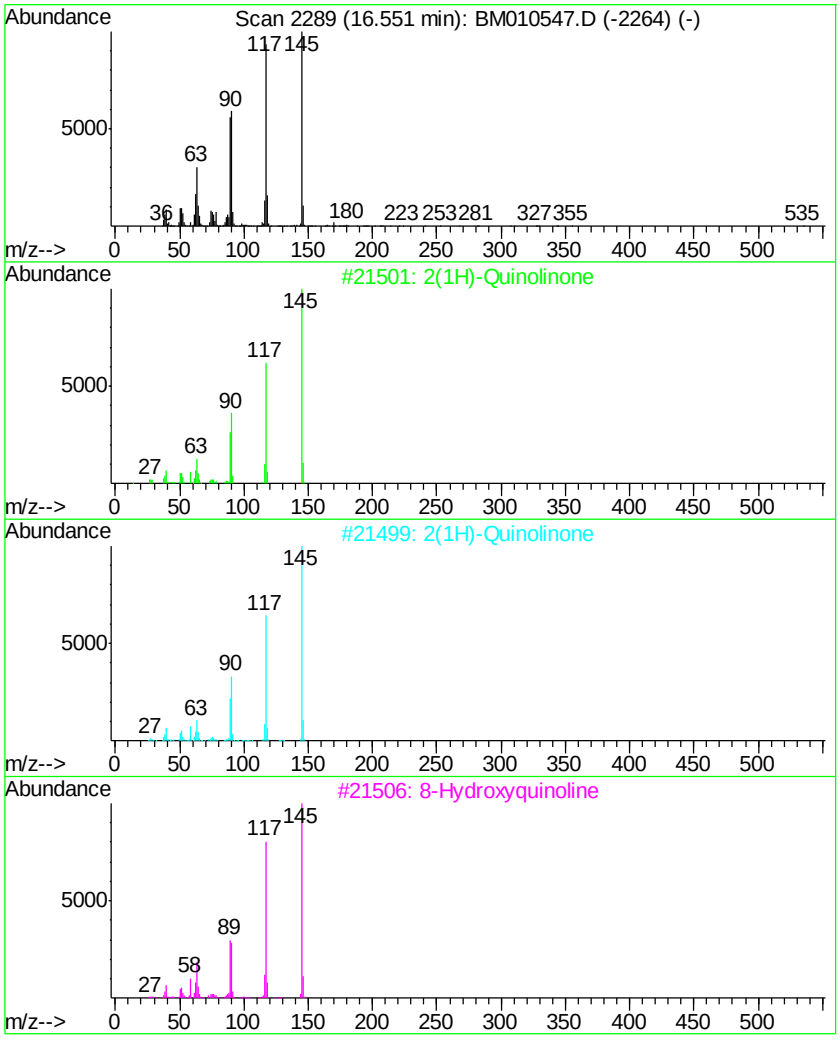
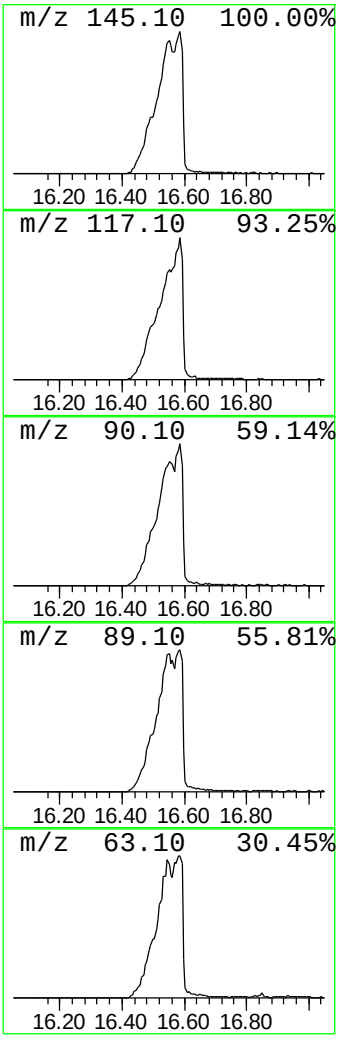
Peak Number 10 2(1H)-Quinolinone Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.55	92.00 ng/ul	9630250	Phenanthrene-d10	17.27

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2(1H)-Quinolinone	145	C9H7NO	000059-31-4	96
2	2(1H)-Quinolinone	145	C9H7NO	000059-31-4	96
3	8-Hydroxyquinoline	145	C9H7NO	000148-24-3	96
4	5-Isoquinolinol	145	C9H7NO	002439-04-5	94
5	5-Quinolinol	145	C9H7NO	000578-67-6	94





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Operator : SJ/MA
Sample : I3558-14DL 2X
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : F5C87DL

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM052717.M
Quant Title : SVOA CALIBRATION

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TIC Integration Parameters: LSCINT.P

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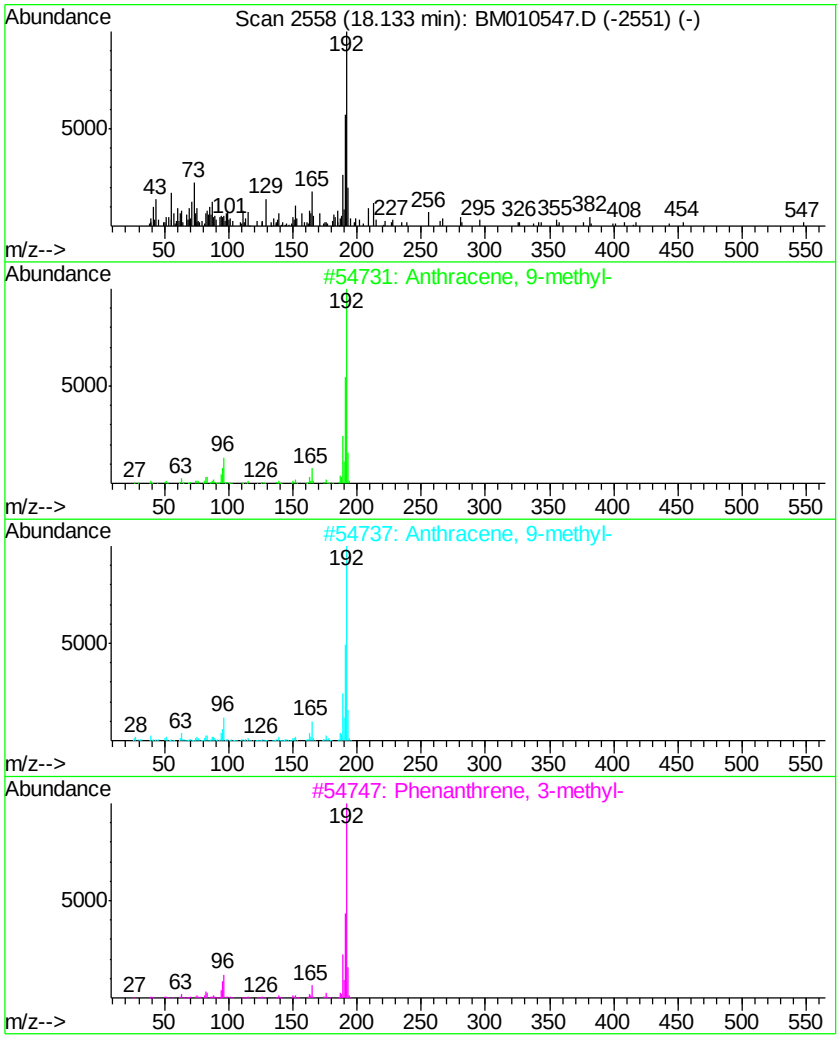
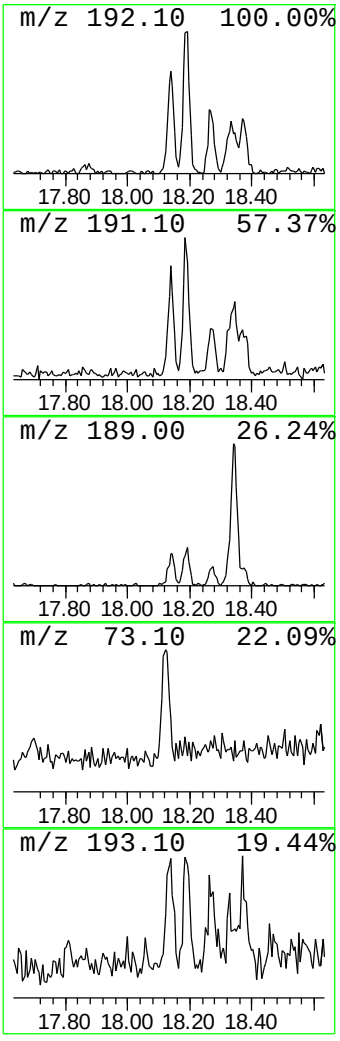
Peak Number 11 unknown-03 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.13	3.56 ng/ul	372364	Phenanthrene-d10	17.27

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Anthracene, 9-methyl-	192	C15H12	000779-02-2	38
2	Anthracene, 9-methyl-	192	C15H12	000779-02-2	38
3	Phenanthrene, 3-methyl-	192	C15H12	000832-71-3	30
4	Anthracene, 1-methyl-	192	C15H12	000610-48-0	30
5	Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	30





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Data File : BM010547.D
Acq On : 12 Jun 2017 22:21
Operator : SJ/MA
Sample : I3558-14DL 2X
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : F5C87DL

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM052717.M
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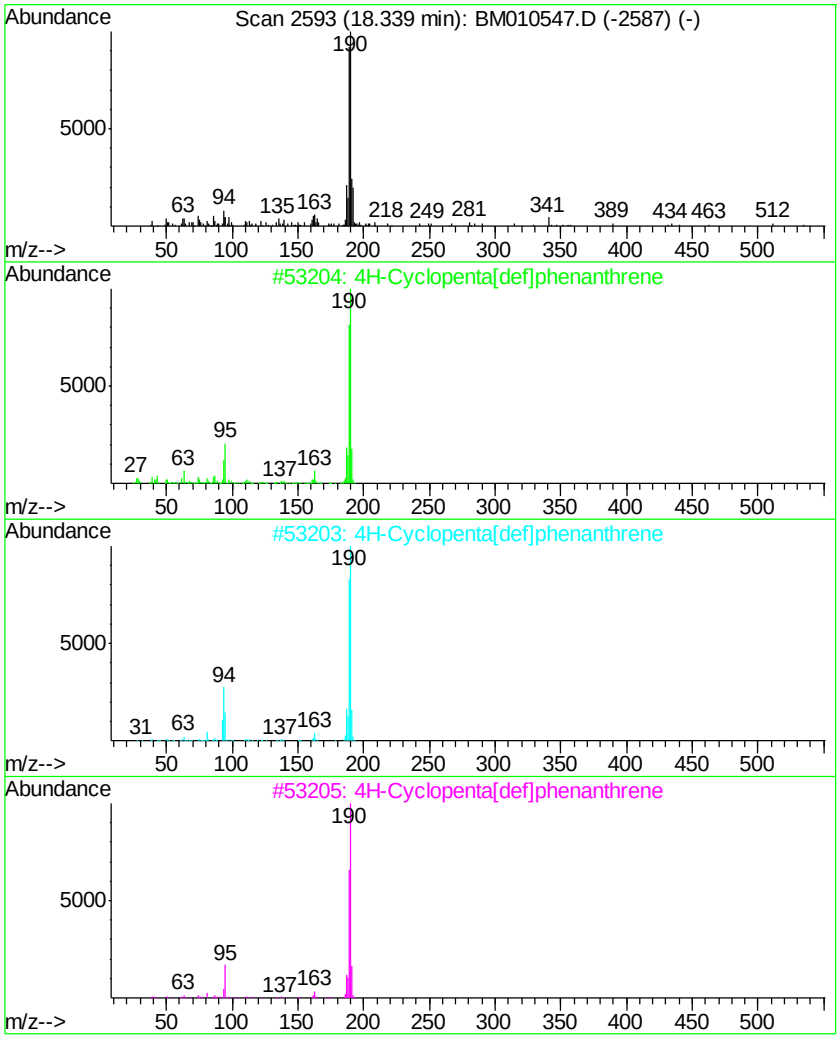
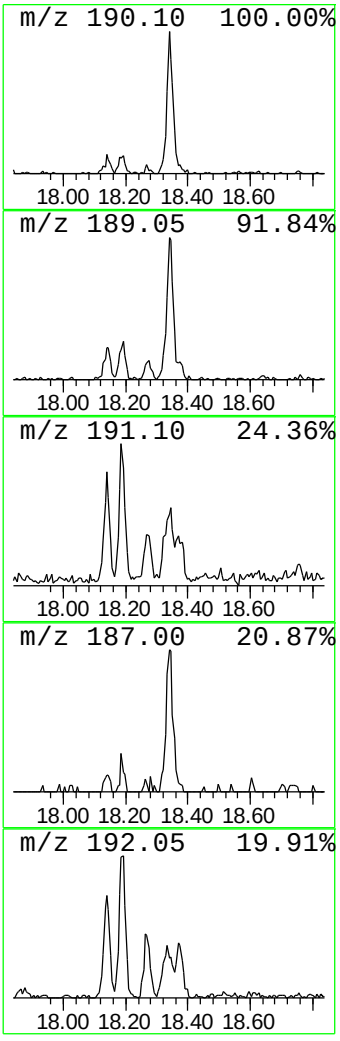
Peak Number 12 4H-Cyclopenta[def]phenanthrene Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.34	2.45 ng/ul	256351	Phenanthrene-d10	17.27

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	81
2	4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	74
3	4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	72
4	Methyl 3beta-acetoxylup-20(29)-e...	512	C33H52O4	004356-30-3	47
5	2,2'-Bis(4,5-dimethylimidazole)	190	C10H14N4	069286-06-2	42





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Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : F5C87DL

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM052717.M
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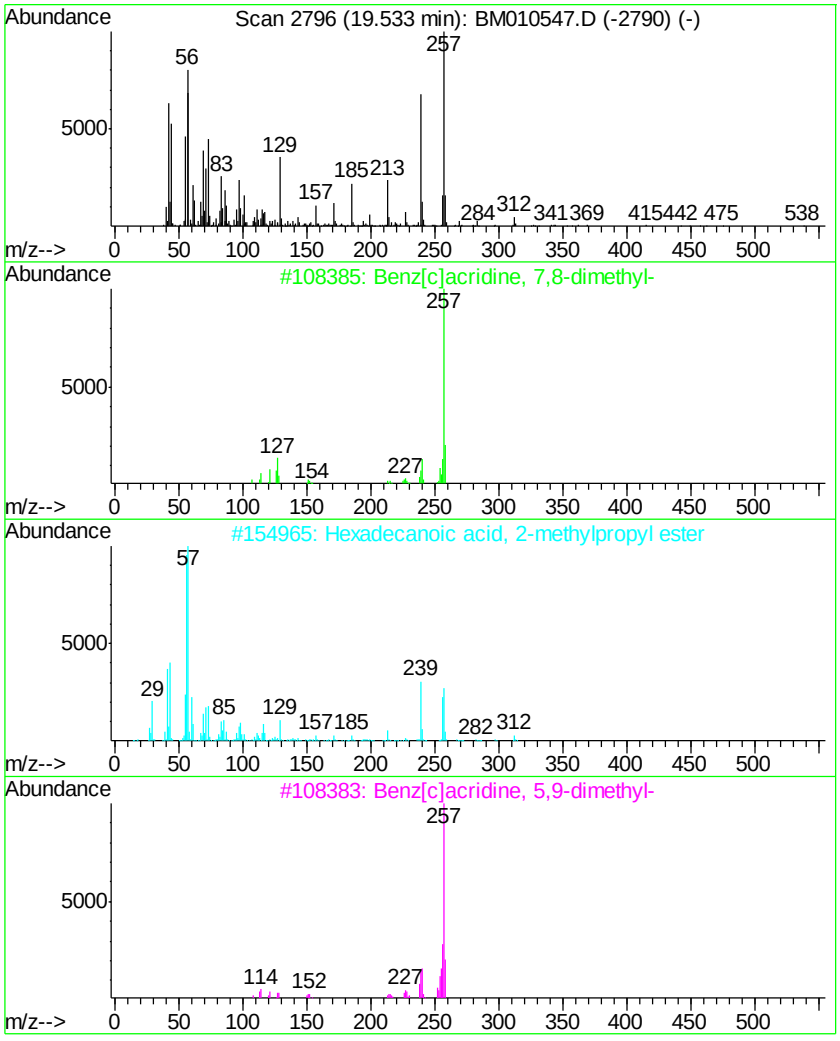
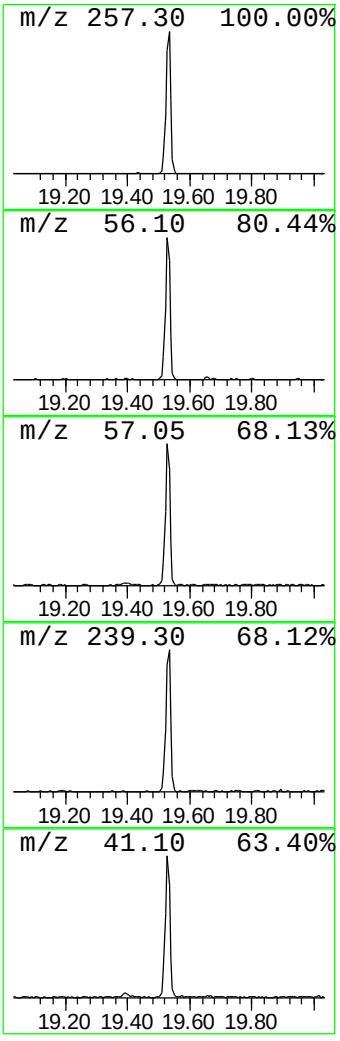
Peak Number 13 unknown-04 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.53	7.93 ng/ul	1384260	Chrysene-d12	21.44

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benz[c]acridine, 7,8-dimethyl-	257	C19H15N	003518-01-2	38
2	Hexadecanoic acid, 2-methylpropy...	312	C20H40O2	000110-34-9	38
3	Benz[c]acridine, 5,9-dimethyl-	257	C19H15N	003518-03-4	38
4	Butyl 4,8,12-trimethyl-tridecanoate	312	C20H40O2	1000336-63-2	38
5	Hexadecanoic acid, butyl ester	312	C20H40O2	000111-06-8	38





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Acq On : 12 Jun 2017 22:21
Operator : SJ/MA
Sample : I3558-14DL 2X
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : F5C87DL

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM052717.M
Quant Title : SVOA CALIBRATION

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

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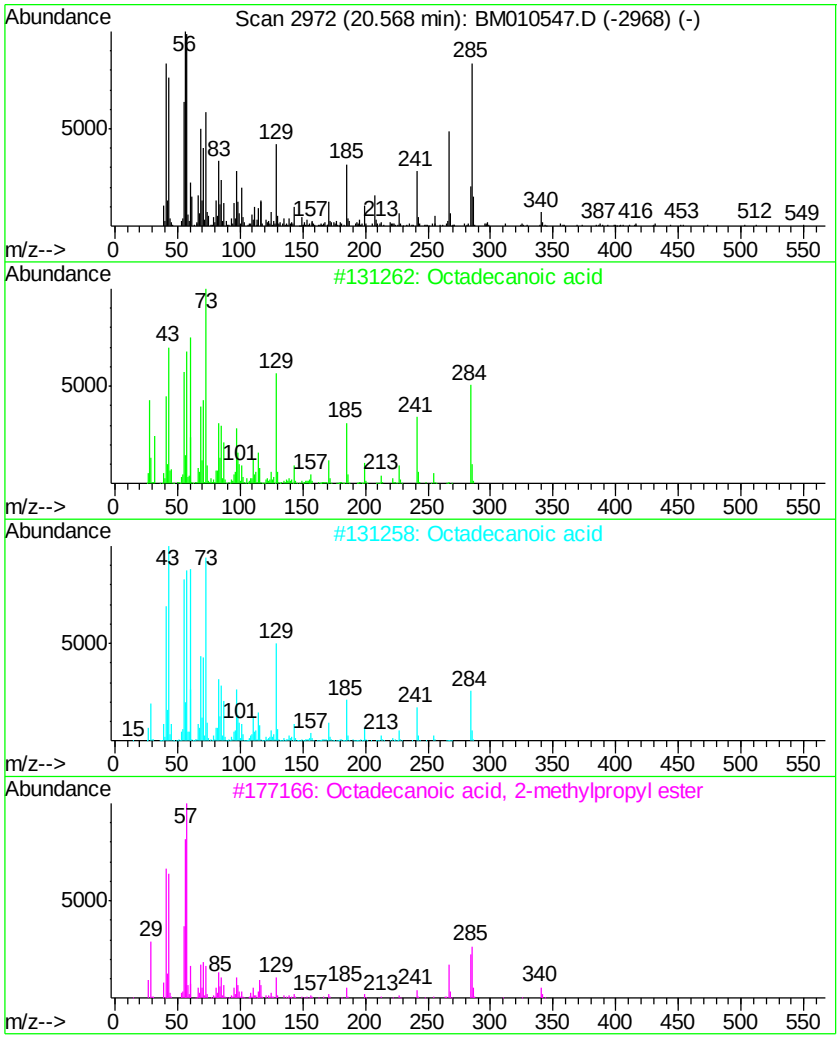
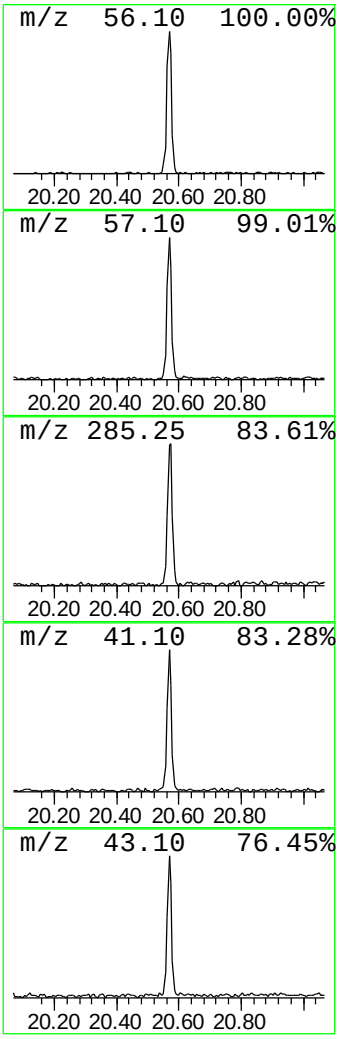
Peak Number 14 Octadecanoic acid Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.57	4.71 ng/ul	821239	Chrysene-d12	21.44

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octadecanoic acid	284	C18H36O2	000057-11-4	62
2	Octadecanoic acid	284	C18H36O2	000057-11-4	53
3	Octadecanoic acid, 2-methylpropy...	340	C22H44O2	000646-13-9	46
4	4',7-Difluoro-2-phenylcinchonini...	285	C16H9F2NO2	020843-19-0	20
5	Benz[c]acridine, 9,11-diethyl-	285	C21H19N	003518-06-7	16





Tentatively Identified Compound (LSC) summary

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM061217\
Data File : BM010547.D
Acq On : 12 Jun 2017 22:21
Operator : SJ/MA
Sample : I3558-14DL 2X
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : F5C87DL

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM052717.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
unknown-01	4.01	5.0	ng/ul	205320	1	7.89	815728	20.0
Phenol, 3,5-dimet...	10.17	3.1	ng/ul	160464	2	10.69	1041370	20.0
Benzoic acid, 3-m...	11.25	4.0	ng/ul	205900	2	10.69	1041370	20.0
Phenol, 2-propyl-	11.52	2.5	ng/ul	130572	2	10.69	1041370	20.0
Cyclohexanemethan...	12.42	12.0	ng/ul	624171	2	10.69	1041370	20.0
Naphthalene, 1-me...	12.56	5.3	ng/ul	276207	2	10.69	1041370	20.0
unknown-02	12.70	3.0	ng/ul	227101	3	14.52	1491020	20.0
2H-Indol-2-one, 1...	14.37	3.9	ng/ul	293751	3	14.52	1491020	20.0
2(1H)-Quinolinone	16.55	92.0	ng/ul	9630250	4	17.27	2093620	20.0
unknown-03	18.13	3.6	ng/ul	372364	4	17.27	2093620	20.0
4H-Cyclopenta[def...	18.34	2.5	ng/ul	256351	4	17.27	2093620	20.0
unknown-04	19.53	7.9	ng/ul	1384260	5	21.44	3490000	20.0
Octadecanoic acid	20.57	4.7	ng/ul	821239	5	21.44	3490000	20.0