LSC Area Percent Report

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM062515\
Data File : BM001796.D
Acq On : 25 Jun 2015 13:35
Operator : TP/IZ
Sample : G2762-03
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : C01L2

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:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM061515.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	2.810	17	21	28	rVB2	18888	29108	0.25%	0.099%
2	2.881	29	33	45	rVB	148450	174360	1.52%	0.591%
3	3.405	119	122	130	rBV	9937	13305	0.12%	0.045%
4	6.175	589	593	601	rBB	14273	21347	0.19%	0.072%
5	6.828	698	704	714	rBB2	131989	279772	2.44%	0.948%
6	6.998	728	733	739	rBB	100239	145291	1.27%	0.492%
7	7.192	760	766	772	rBB	132638	199582	1.74%	0.676%
8	7.663	840	846	852	rBB	168261	259782	2.27%	0.880%
9	8.098	916	920	926	rBB	39352	57070	0.50%	0.193%
10	8.363	958	965	970	rBV	164357	257183	2.24%	0.871%
11	8.434	970	977	984	rVB	329433	575441	5.02%	1.949%
12	8.822	1037	1043	1049	rBB	118680	188737	1.65%	0.639%
13	9.422	1141	1145	1150	rBB	16065	21946	0.19%	0.074%
14	9.539	1159	1165	1171	rBB	108039	168653	1.47%	0.571%
15	9.628	1174	1180	1183	rBV	205119	328627	2.87%	1.113%
16	9.663	1183	1186	1192	rVB	147173	209585	1.83%	0.710%
17	9.892	1217	1225	1229	rBV2	78684	189278	1.65%	0.641%
18	9.934	1229	1232	1239	rVB	99410	143077	1.25%	0.485%
19	10.075	1249	1256	1264	rBB2	191139	389550	3.40%	1.320%
20	10.322	1293	1298	1304	rBV	87324	126826	1.11%	0.430%
21	10.451	1313	1320	1327	rBB	221430	358855	3.13%	1.216%
22	10.592	1337	1344	1351	rBB	171157	269201	2.35%	0.912%
23	10.804	1375	1380	1386	rBB	14827	26179	0.23%	0.089%
24	10.986	1406	1411	1416	rBV	24065	35225	0.31%	0.119%
25	11.069	1420	1425	1427	rVV	14485	23001	0.20%	0.078%
26	11.098	1427	1430	1439	rVB	21947	36065	0.31%	0.122%
27	11.275	1455	1460	1466	rBV2	43469	69884	0.61%	0.237%
28	11.475	1488	1494	1499	rBV2	12841	18335	0.16%	0.062%
29	11.528	1499	1503	1507	rVV2	10204	15737	0.14%	0.053%
30	11.575	1507	1511	1516	rVB2	13707	21549	0.19%	0.073%
31	13.727	1871	1877	1881	rBV	284948	371972	3.25%	1.260%
32	13.774	1881	1885	1890	rVB	110891	145011	1.27%	0.491%
33	14.010	1918	1925	1932	rVB	292400	439883	3.84%	1.490%
34	14.316	1971	1977	1983	rBV2	332327	489682	4.27%	1.659%

LSC Area Percent Report

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM062515\
Data File : BM001796.D
Acq On : 25 Jun 2015 13:35
Operator : TP/IZ
Sample : G2762-03
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : C01L2

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:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM061515.M
Title : SVOA CALIBRATION

35	14.516	2005	2011	2020	rBV	140956	198886	1.74%	0.674%
36	15.310	2140	2146	2152	rVB	396284	551255	4.81%	1.867%
37	15.433	2161	2167	2174	rVB2	111203	151593	1.32%	0.514%
38	15.527	2177	2183	2188	rBV	27255	32597	0.28%	0.110%
39	16.068	2258	2275	2280	rBV	1941801	5739248	50.07%	19.441%
40	16.421	2330	2335	2341	rVB	55646	74647	0.65%	0.253%
41	17.062	2438	2444	2450	rBV	454931	605569	5.28%	2.051%
42	17.162	2452	2461	2467	rBV	462965	632665	5.52%	2.143%
43	17.980	2595	2600	2606	rBV	115391	148155	1.29%	0.502%
44	18.362	2647	2665	2669	rBV	3961368	11462495	100.00%	38.828%
45	18.504	2686	2689	2696	rVB	89285	109766	0.96%	0.372%
46	19.245	2810	2815	2827	rVV2	70313	96427	0.84%	0.327%
47	19.351	2829	2833	2836	rVV2	13764	17782	0.16%	0.060%
48	19.392	2836	2840	2844	rVV4	7318	11634	0.10%	0.039%
49	19.462	2846	2852	2864	rVB	572356	767157	6.69%	2.599%
50	19.792	2904	2908	2914	rBV5	11186	16339	0.14%	0.055%
51	19.898	2922	2926	2935	rBV4	10805	17619	0.15%	0.060%
52	20.080	2953	2957	2961	rBV4	12867	16816	0.15%	0.057%
53	20.150	2965	2969	2974	rVB3	20465	26390	0.23%	0.089%
54	20.239	2974	2984	2991	rBV	76426	103626	0.90%	0.351%
55	20.492	3023	3027	3031	rVV3	9895	16202	0.14%	0.055%
56	20.827	3078	3084	3088	rBV3	9753	14597	0.13%	0.049%
57	20.968	3103	3108	3115	rBV	142676	185151	1.62%	0.627%
58	21.021	3115	3117	3120	rVV2	10291	11548	0.10%	0.039%
59	21.174	3139	3143	3150	rVB5	17158	22222	0.19%	0.075%
60	21.256	3150	3157	3162	rBV	653598	797056	6.95%	2.700%
61	21.850	3253	3258	3261	rBV6	14629	23557	0.21%	0.080%
62	23.121	3469	3474	3480	rVB3	20492	32503	0.28%	0.110%
63	23.333	3503	3510	3521	rBV2	419212	769555	6.71%	2.607%
64	23.480	3527	3535	3544	rVB	387990	713175	6.22%	2.416%
65	24.080	3630	3637	3646	rBV7	21640	55592	0.48%	0.188%

Sum of corrected areas: 29520923

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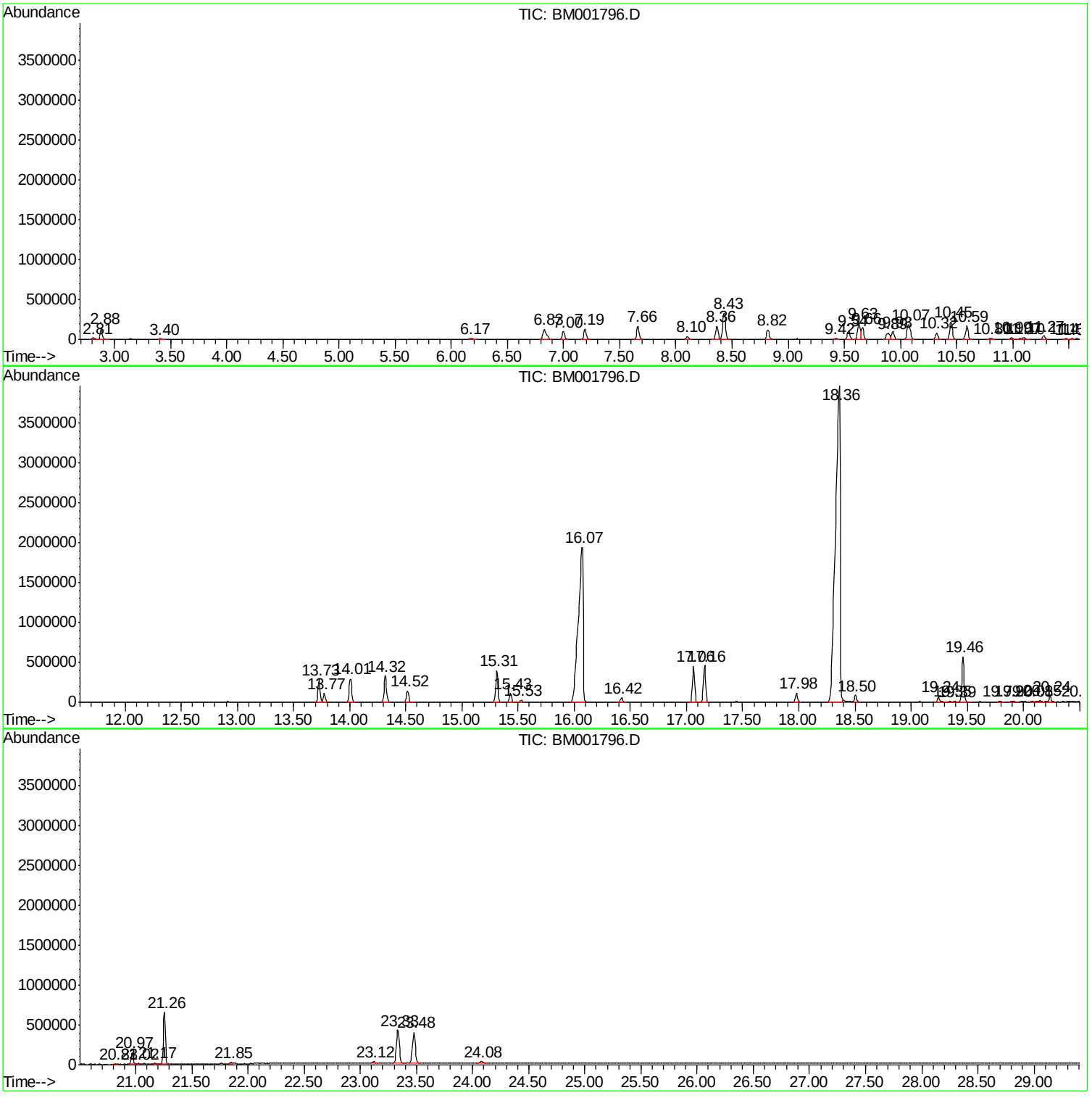
Acq On : 25 Jun 2015 13:35
Operator : TP/IZ
Sample : G2762-03
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
C01L2

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM061515.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P



Library Search Compound Report

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM062515\
Data File : BM001796.D
Acq On : 25 Jun 2015 13:35
Operator : TP/IZ
Sample : G2762-03
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : C01L2

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM061515.M
Quant Title : SVOA CALIBRATION

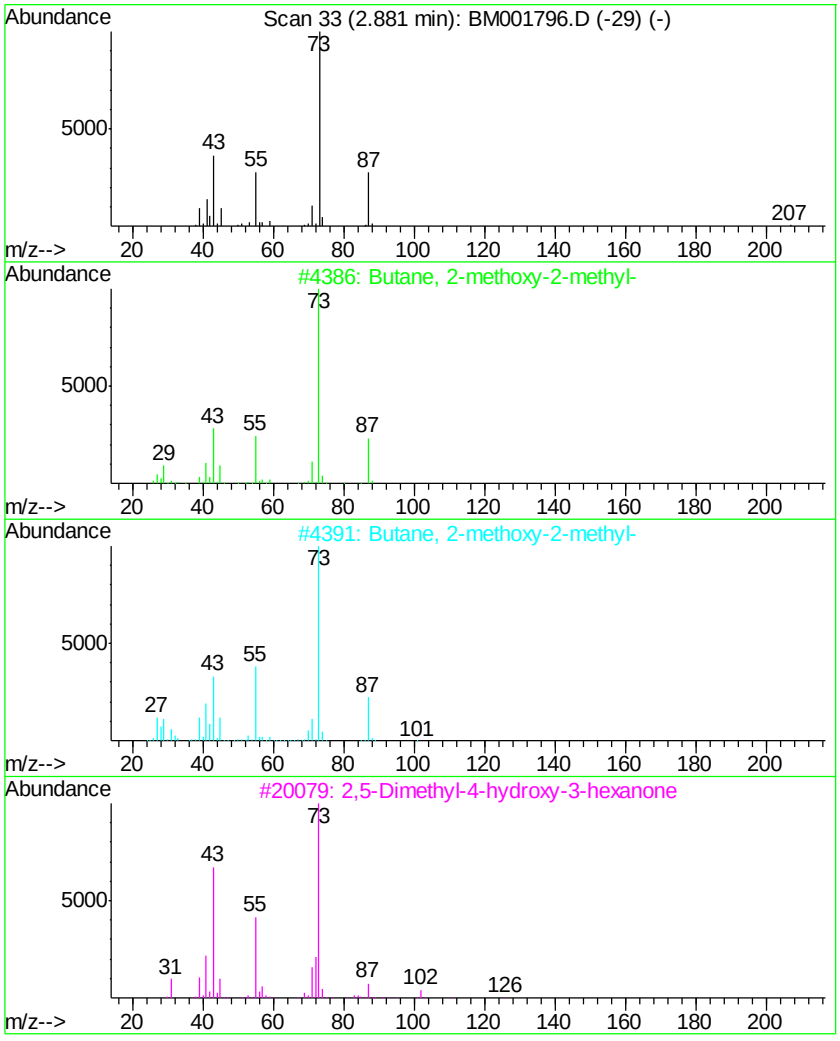
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

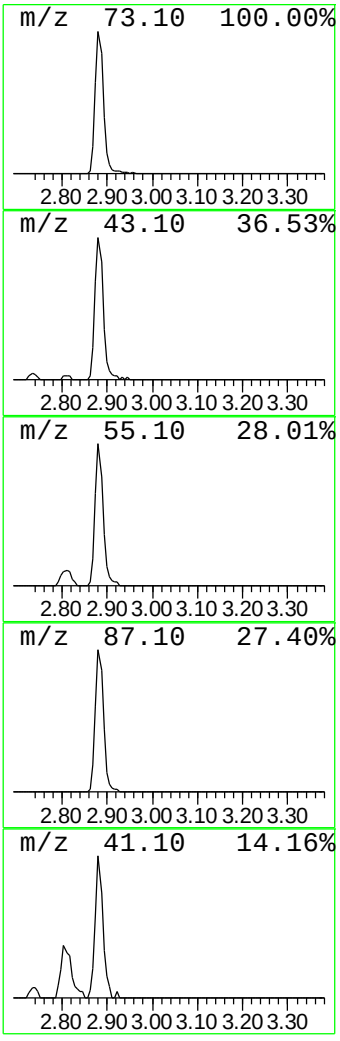
****

Peak Number 2 Butane, 2-methoxy-2-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.88	13.42 ng/ul	174360	1,4-Dichlorobenzene-d4	7.66

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	78
2	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	74
3	2,5-Dimethyl-4-hydroxy-3-hexanone	144	C8H16O2	000815-77-0	25
4	Silane, tetramethyl-	88	C4H12Si	000075-76-3	23
5	Pentane, 3-methoxy-	102	C6H14O	036839-67-5	17





Library Search Compound Report

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM062515\
Data File : BM001796.D
Acq On : 25 Jun 2015 13:35
Operator : TP/IZ
Sample : G2762-03
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : C01L2

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM061515.M
Quant Title : SVOA CALIBRATION

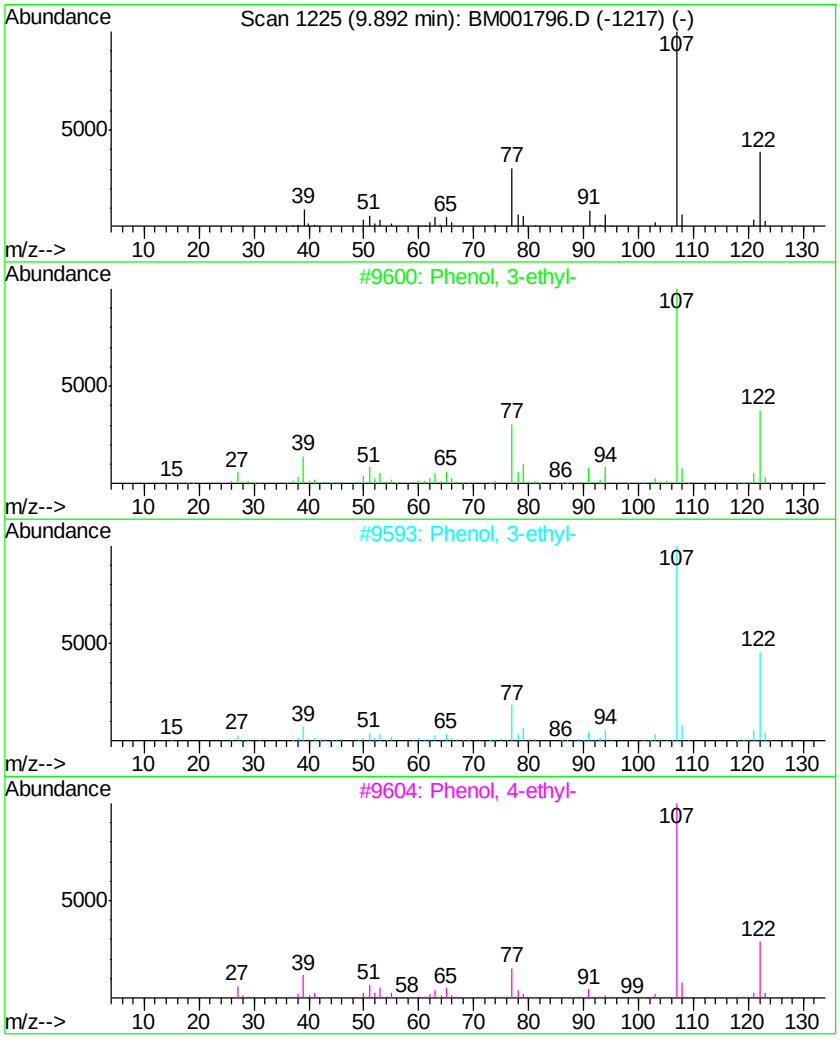
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

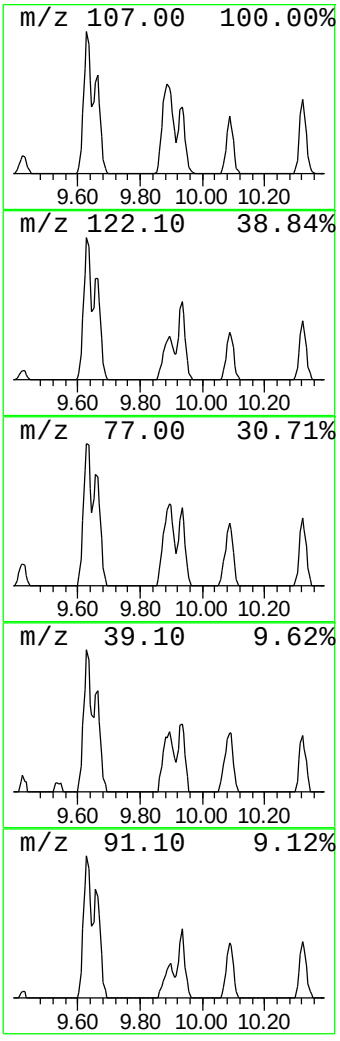
****

Peak Number 3 Phenol, 3-ethyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.89	10.55 ng/ul	189278	Naphthalene-d8	10.45

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, 3-ethyl-	122	C8H10O	000620-17-7	97
2	Phenol, 3-ethyl-	122	C8H10O	000620-17-7	91
3	Phenol, 4-ethyl-	122	C8H10O	000123-07-9	90
4	Phenol, 2-ethyl-	122	C8H10O	000090-00-6	90
5	Phenol, 2-ethyl-	122	C8H10O	000090-00-6	87





Data Path : Z:\HPCHEM1\BNA_M\DATA\BM062515\
Data File : BM001796.D
Acq On : 25 Jun 2015 13:35
Operator : TP/IZ
Sample : G2762-03
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : C01L2

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM061515.M
Quant Title : SVOA CALIBRATION

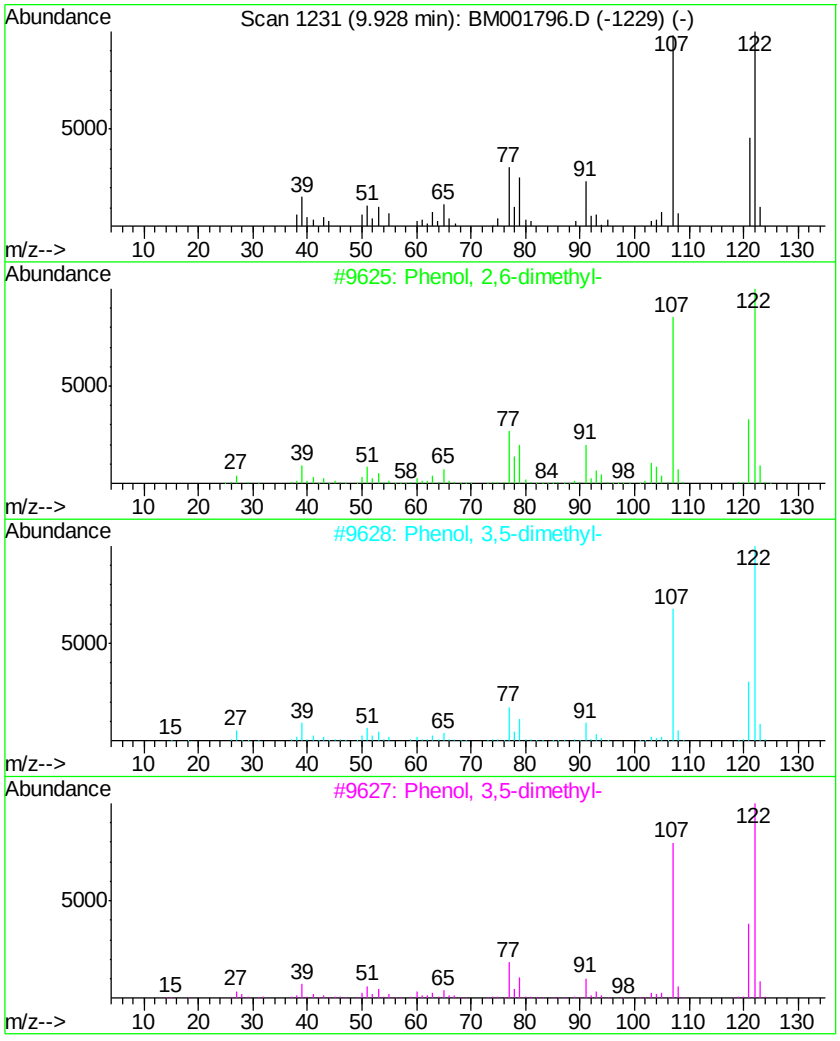
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

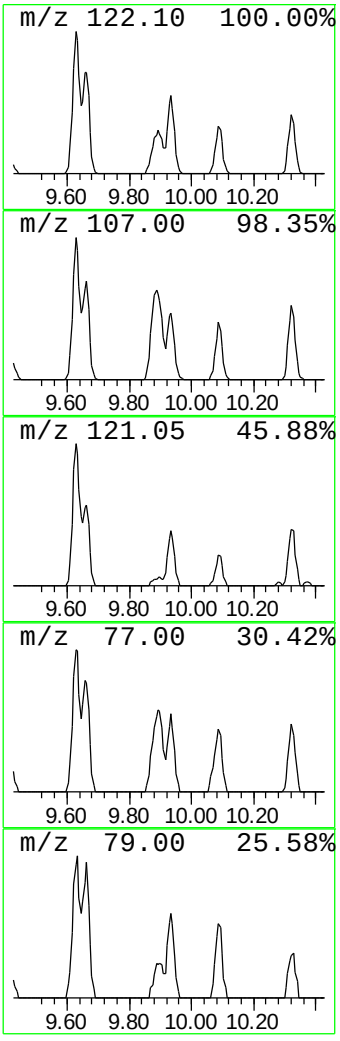
****

Peak Number 4 Phenol, 2,6-dimethyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.93	7.97 ng/ul	143077	Naphthalene-d8	10.45

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, 2,6-dimethyl-	122	C8H10O	000576-26-1	91
2	Phenol, 3,5-dimethyl-	122	C8H10O	000108-68-9	91
3	Phenol, 3,5-dimethyl-	122	C8H10O	000108-68-9	91
4	Phenol, 2,3-dimethyl-	122	C8H10O	000526-75-0	91
5	Phenol, 2,3-dimethyl-	122	C8H10O	000526-75-0	91





Data Path : Z:\HPCHEM1\BNA_M\DATA\BM062515\
Data File : BM001796.D
Acq On : 25 Jun 2015 13:35
Operator : TP/IZ
Sample : G2762-03
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : C01L2

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM061515.M
Quant Title : SVOA CALIBRATION

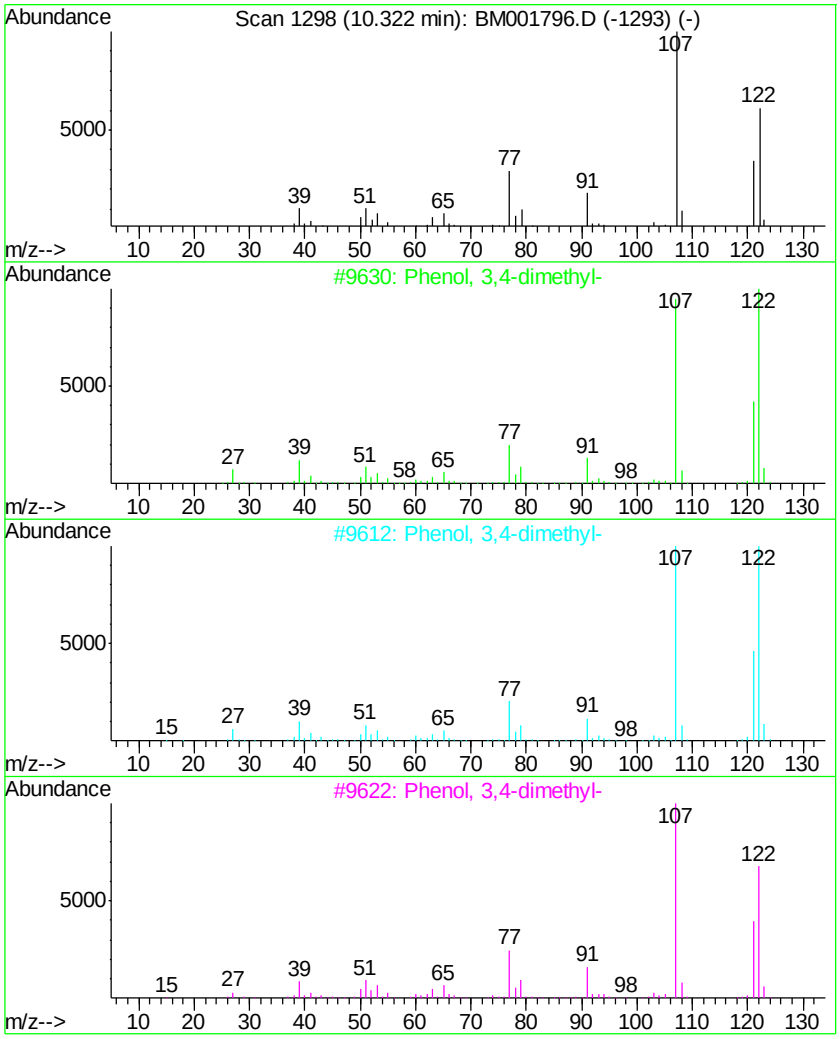
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

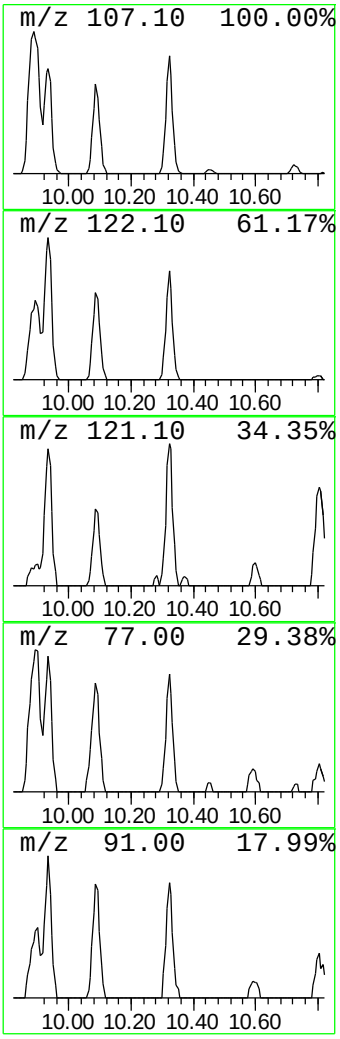
****

Peak Number 5 Phenol, 3,4-dimethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.32	7.07 ng/ul	126826	Naphthalene-d8	10.45

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, 3,4-dimethyl-	122	C8H10O	000095-65-8	94
2	Phenol, 3,4-dimethyl-	122	C8H10O	000095-65-8	94
3	Phenol, 3,4-dimethyl-	122	C8H10O	000095-65-8	91
4	Phenol, 3-ethyl-	122	C8H10O	000620-17-7	91
5	Phenol, 3-ethyl-	122	C8H10O	000620-17-7	91





Library Search Compound Report

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM062515\
Data File : BM001796.D
Acq On : 25 Jun 2015 13:35
Operator : TP/IZ
Sample : G2762-03
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : C01L2

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM061515.M
Quant Title : SVOA CALIBRATION

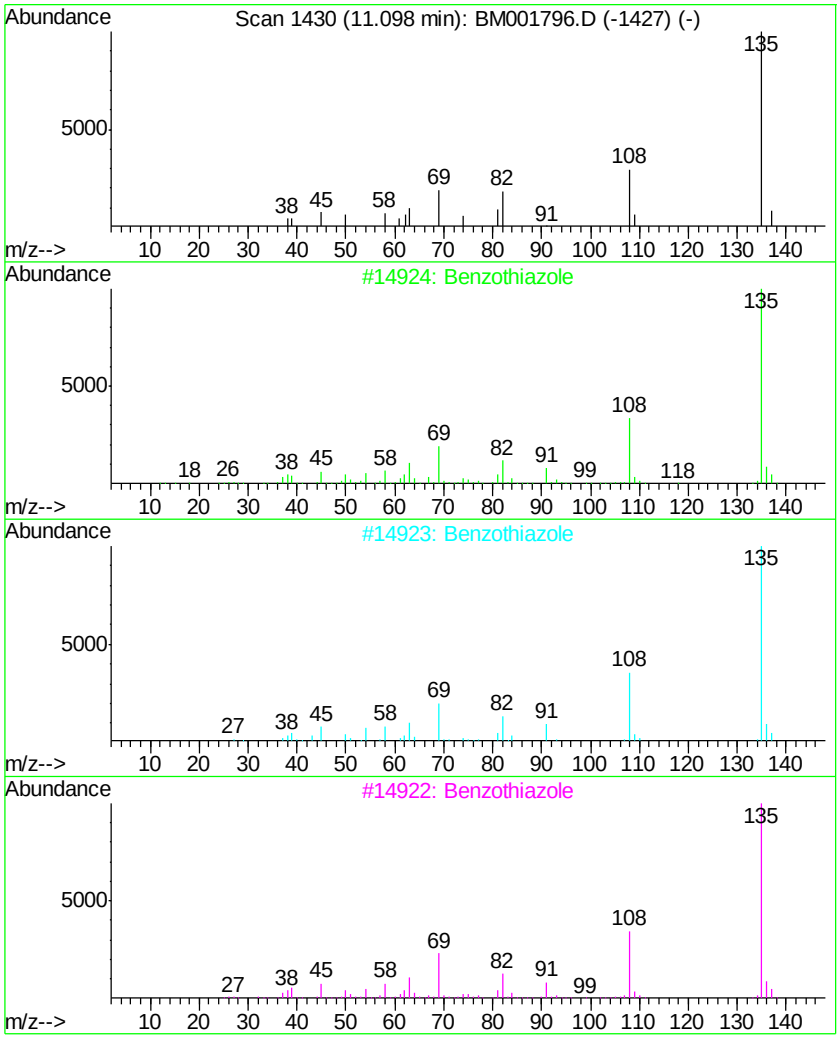
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

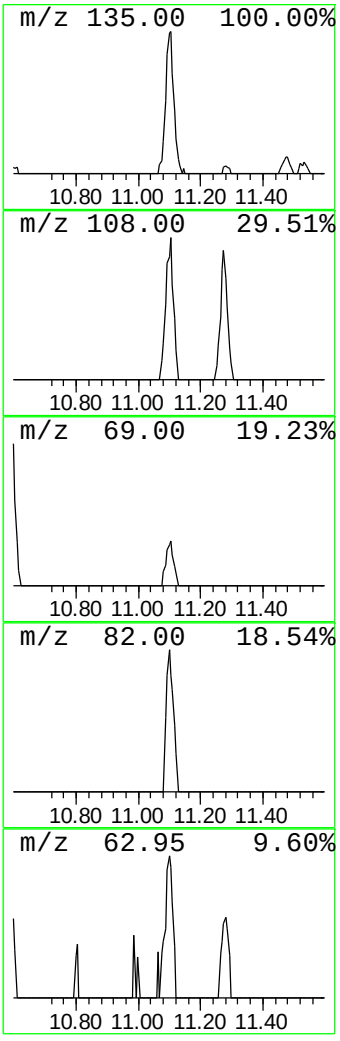
****

Peak Number 6 Benzothiazole Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.10	2.01 ng/ul	36065	Naphthalene-d8	10.45

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzothiazole	135	C7H5NS	000095-16-9	86
2	Benzothiazole	135	C7H5NS	000095-16-9	78
3	Benzothiazole	135	C7H5NS	000095-16-9	78
4	1,2-Benzisothiazole	135	C7H5NS	000272-16-2	72
5	Thieno[2,3-c]pyridine	135	C7H5NS	000272-12-8	72





Data Path : Z:\HPCHEM1\BNA_M\DATA\BM062515\
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Acq On : 25 Jun 2015 13:35
Operator : TP/IZ
Sample : G2762-03
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : C01L2

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM061515.M
Quant Title : SVOA CALIBRATION

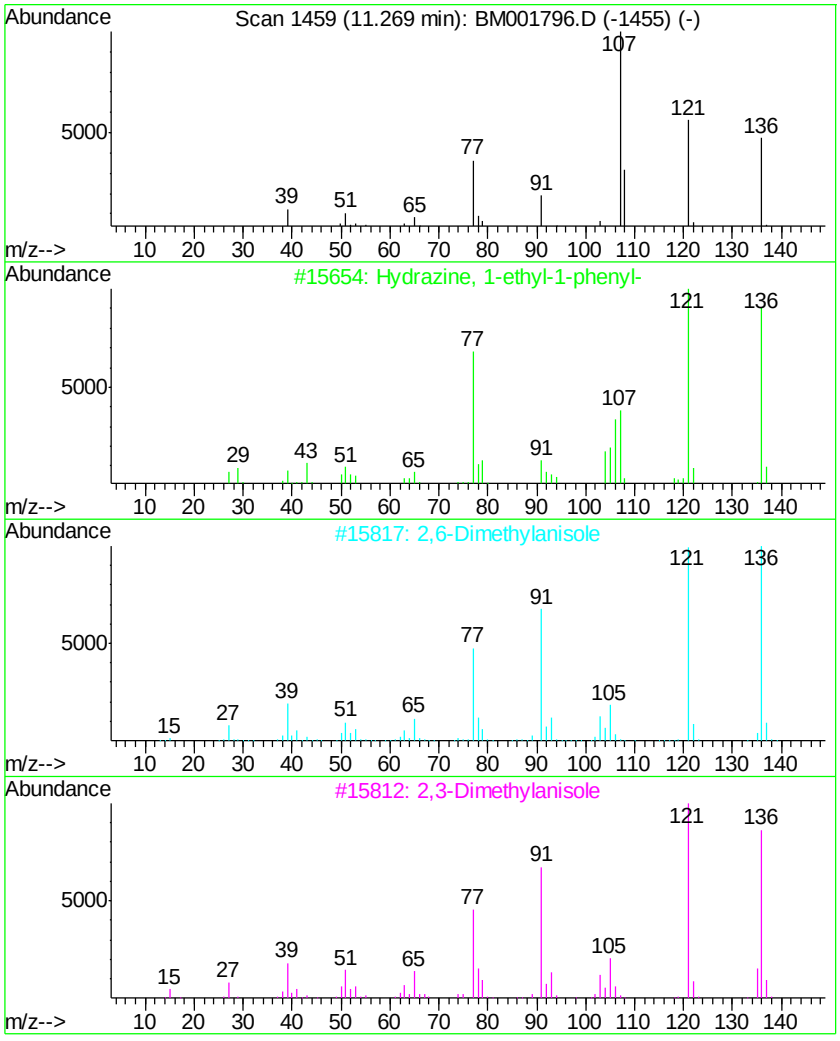
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

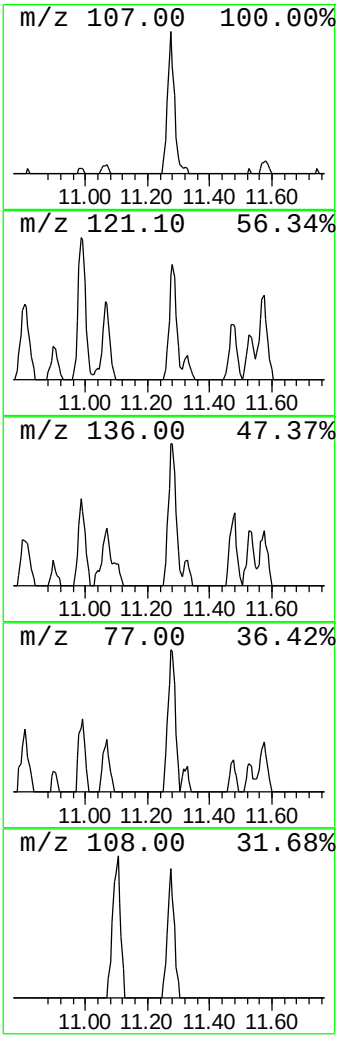
****

Peak Number 7 Hydrazine, 1-ethyl-1-phenyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.27	3.89 ng/ul	69884	Naphthalene-d8	10.45

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hydrazine, 1-ethyl-1-phenyl-	136	C8H12N2	000644-21-3	52
2	2,6-Dimethylanisole	136	C9H12O	001004-66-6	49
3	2,3-Dimethylanisole	136	C9H12O	002944-49-2	49
4	Phenol, 2-propyl-	136	C9H12O	000644-35-9	49
5	Furan, 2-(1-pentenyl)-, (E)-	136	C9H12O	020992-69-2	49





Data Path : Z:\HPCHEM1\BNA_M\DATA\BM062515\
Data File : BM001796.D
Acq On : 25 Jun 2015 13:35
Operator : TP/IZ
Sample : G2762-03
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : C01L2

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM061515.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

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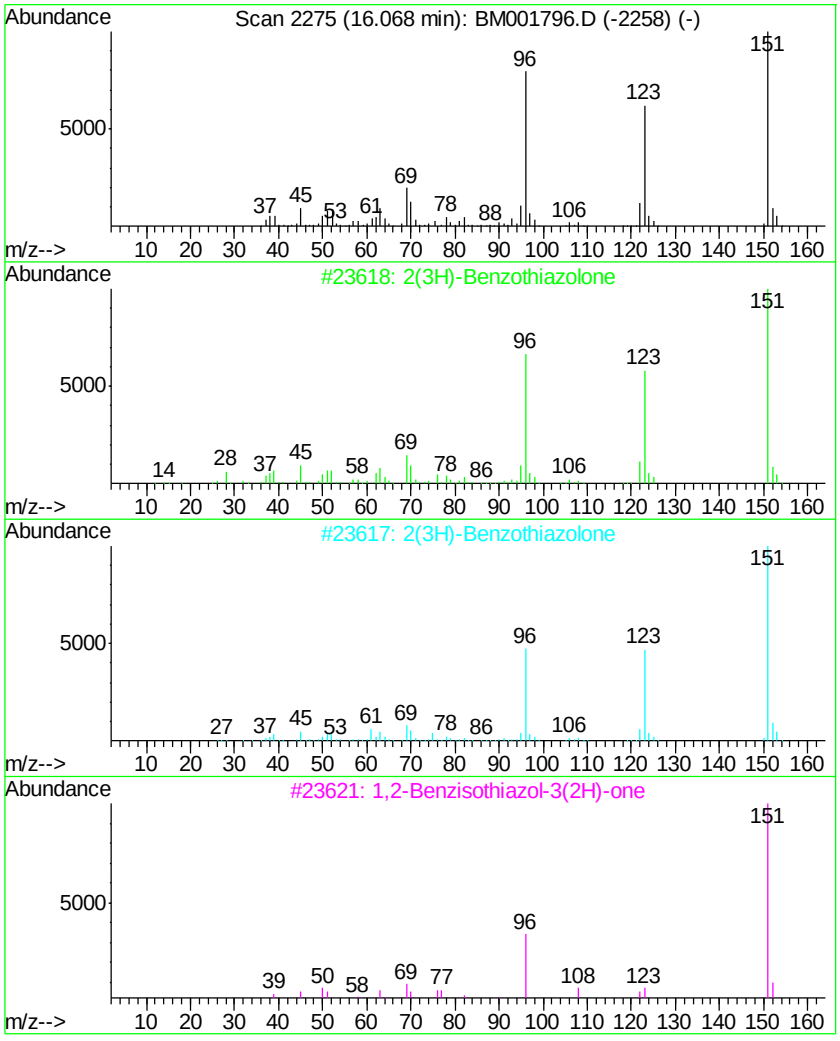
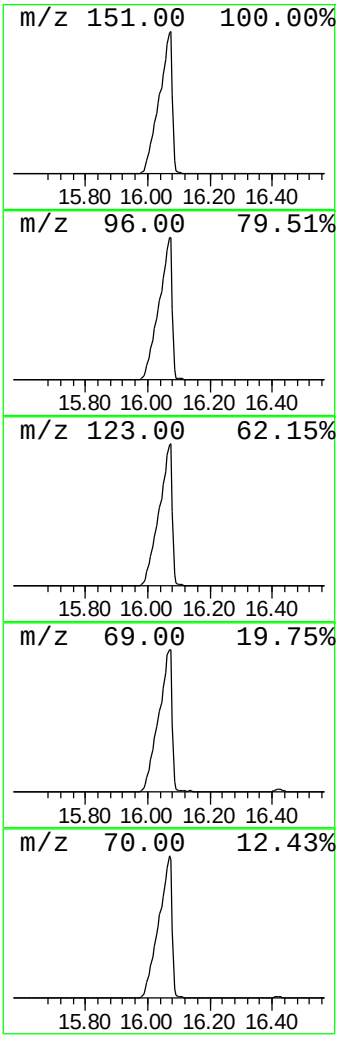
Peak Number 8 2(3H)-Benzothiazolone Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.07	189.55 ng/ul	5739250	Phenanthrene-d10	17.06

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2(3H)-Benzothiazolone	151	C7H5NOS	000934-34-9	97
2	2(3H)-Benzothiazolone	151	C7H5NOS	000934-34-9	94
3	1,2-Benzisothiazol-3(2H)-one	151	C7H5NOS	002634-33-5	59
4	Thieno[2,3-b]pyridine-N-oxide	151	C7H5NOS	025557-50-0	43
5	4-Methoxyformanilide	151	C8H9NO2	023896-88-0	38





Data Path : Z:\HPCHEM1\BNA_M\DATA\BM062515\
Data File : BM001796.D
Acq On : 25 Jun 2015 13:35
Operator : TP/IZ
Sample : G2762-03
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : C01L2

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM061515.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

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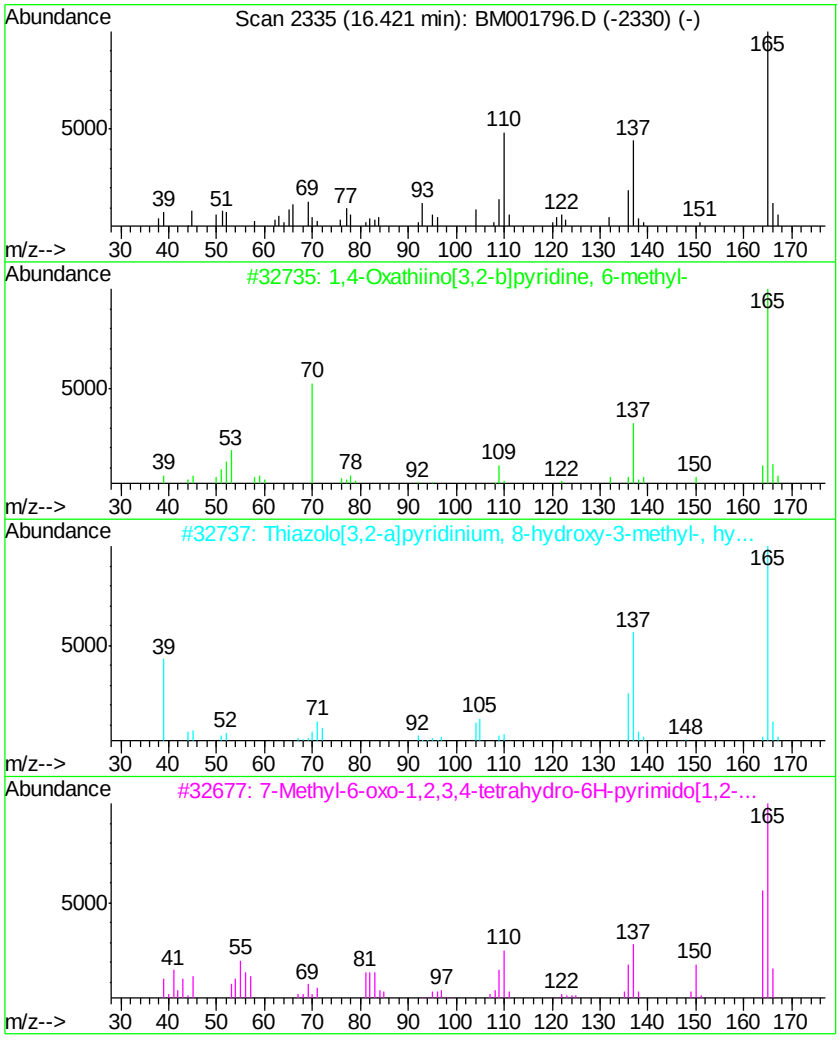
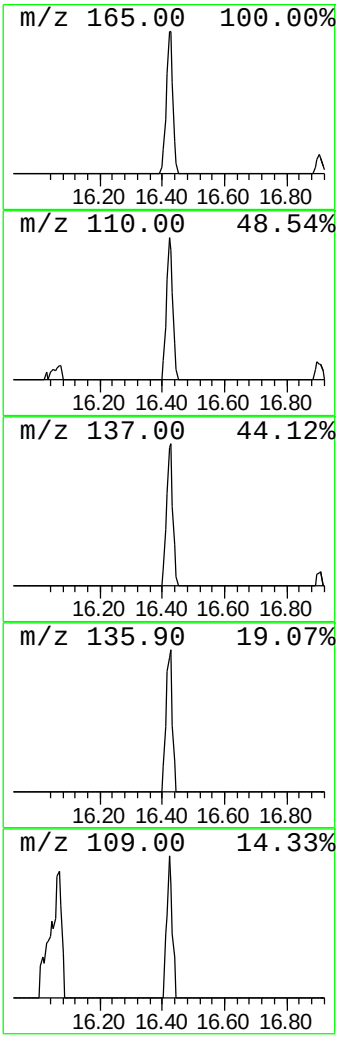
Peak Number 9 1,4-Oxathiino[3,2-b]pyridin... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.42	2.47 ng/ul	74647	Phenanthrene-d10	17.06

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,4-Oxathiino[3,2-b]pyridine, 6-...	165	C8H7NOS	056114-39-7	58
2	Thiazolo[3,2-a]pyridinium, 8-hyd...	165	C8H7NOS	030276-99-4	53
3	7-Methyl-6-oxo-1,2,3,4-tetrahydr...	165	C8H11N3O	026955-14-6	50
4	3,4-Dihydro-benzo[1,4]thiazin-2-one	165	C8H7NOS	096220-47-2	49
5	5-Amino-3-(2-thienyl)pyrazole	165	C7H7N3S	096799-03-0	45





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Data File : BM001796.D
Acq On : 25 Jun 2015 13:35
Operator : TP/IZ
Sample : G2762-03
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : C01L2

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM061515.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

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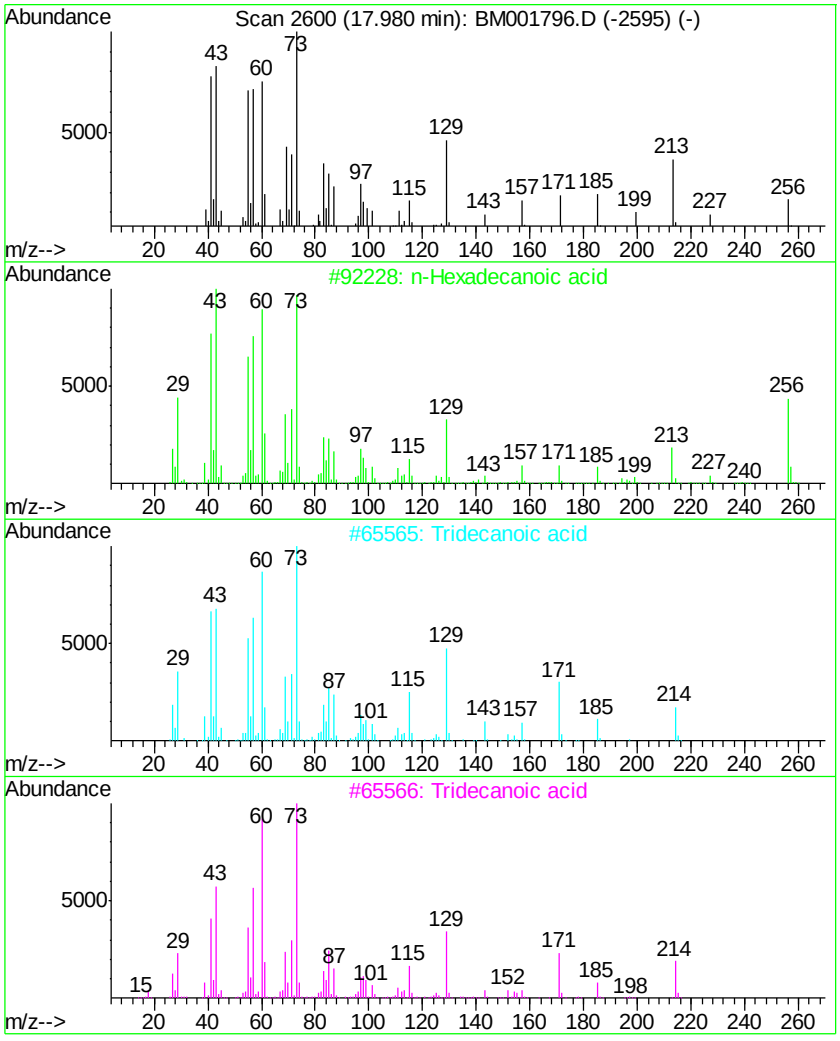
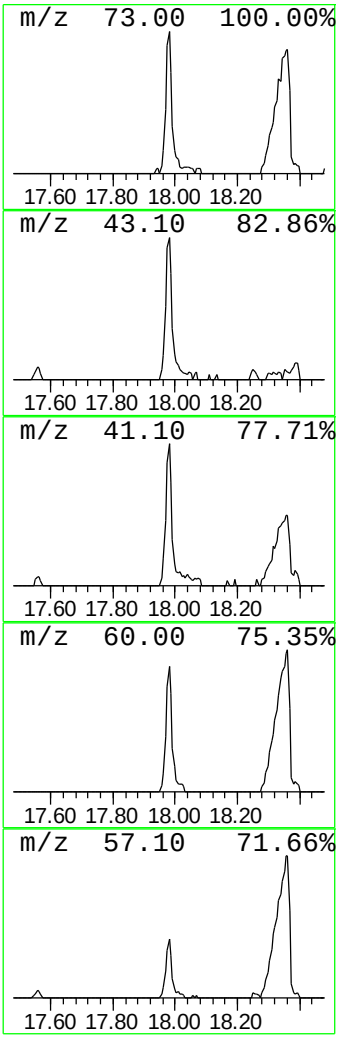
Peak Number 10 n-Hexadecanoic acid Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.98	4.89 ng/ul	148155	Phenanthrene-d10	17.06

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	95
2	Tridecanoic acid	214	C13H26O2	000638-53-9	91
3	Tridecanoic acid	214	C13H26O2	000638-53-9	91
4	Tetradecanoic acid	228	C14H28O2	000544-63-8	80
5	Pentadecanoic acid	242	C15H30O2	001002-84-2	80





Library Search Compound Report

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Data File : BM001796.D
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Operator : TP/IZ
Sample : G2762-03
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : C01L2

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM061515.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

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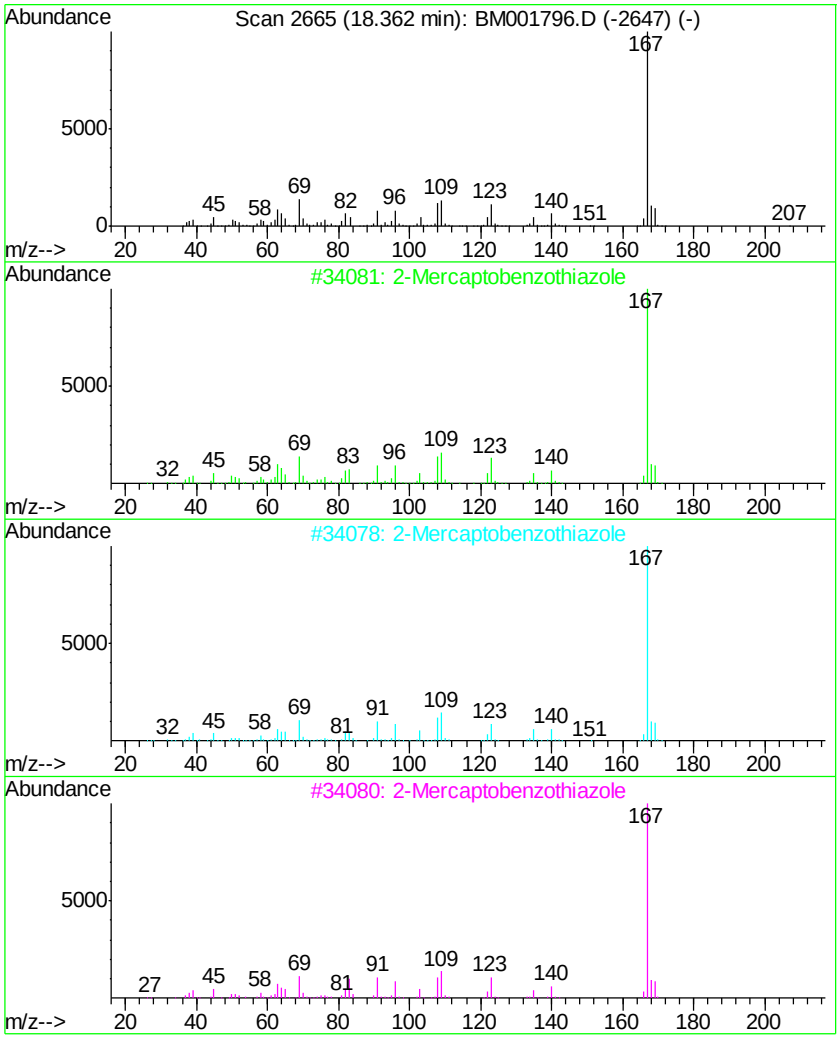
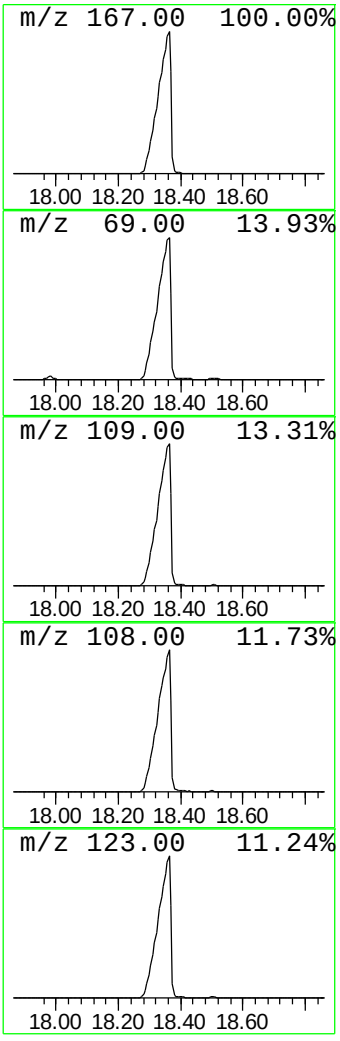
Peak Number 11 2-Mercaptobenzothiazole Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.36	378.57 ng/ul	11462500	Phenanthrene-d10	17.06

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Mercaptobenzothiazole	167	C7H5NS2	000149-30-4	96
2	2-Mercaptobenzothiazole	167	C7H5NS2	000149-30-4	96
3	2-Mercaptobenzothiazole	167	C7H5NS2	000149-30-4	94
4	2-Mercaptobenzothiazole	167	C7H5NS2	000149-30-4	90
5	5-Fluoro-2-methyl-benzothiazole	167	C8H6FNS	000399-75-7	52





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Operator : TP/IZ
Sample : G2762-03
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : C01L2

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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

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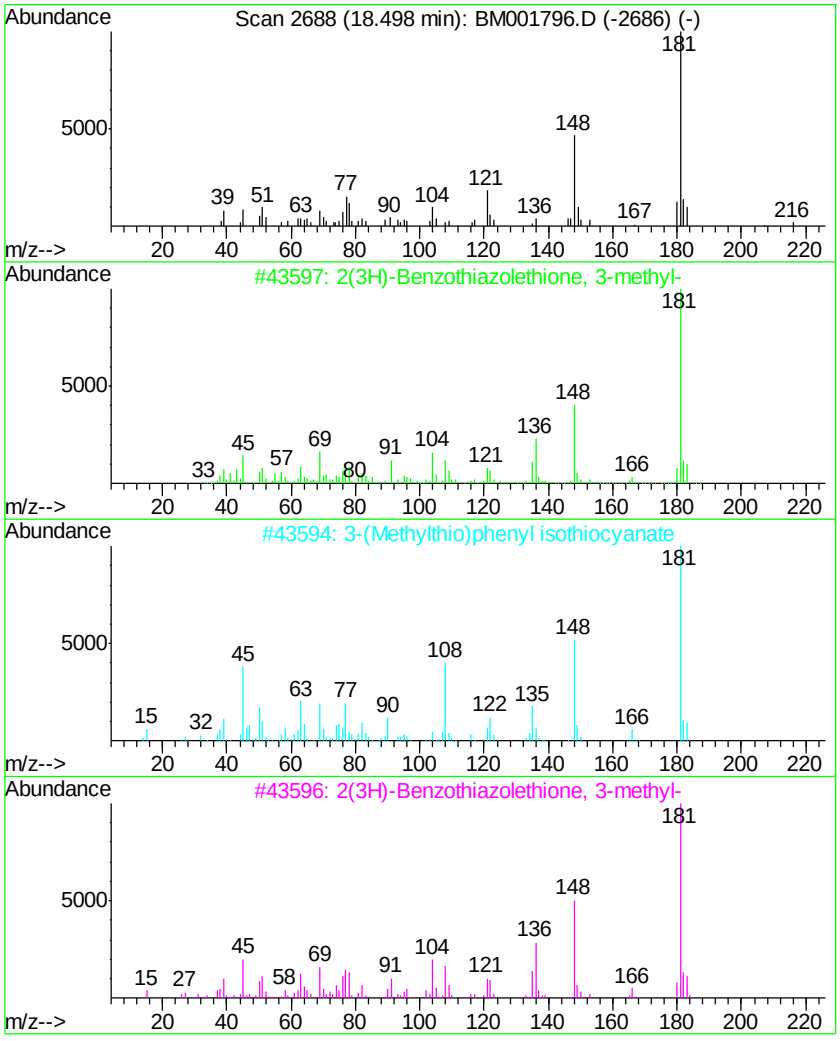
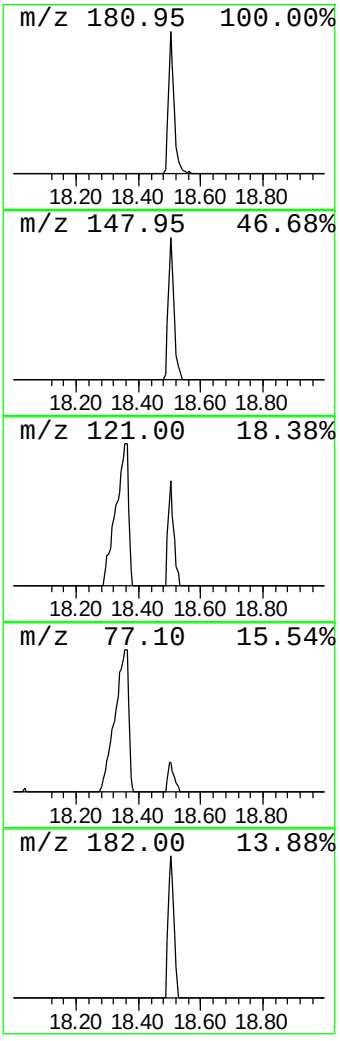
Peak Number 12 2(3H)-Benzothiazolethione, ... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.50	3.63 ng/ul	109766	Phenanthrene-d10	17.06

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2(3H)-Benzothiazolethione, 3-met...	181	C8H7NS2	002254-94-6	87
2	3-(Methylthio)phenyl isothiocyanate	181	C8H7NS2	051333-80-3	72
3	2(3H)-Benzothiazolethione, 3-met...	181	C8H7NS2	002254-94-6	64
4	Benzothiazole, 2-(methylthio)-	181	C8H7NS2	000615-22-5	64
5	Benzothiazole, 2-(methylthio)-	181	C8H7NS2	000615-22-5	64





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Sample : G2762-03
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ALS Vial : 7 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : C01L2

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM061515.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

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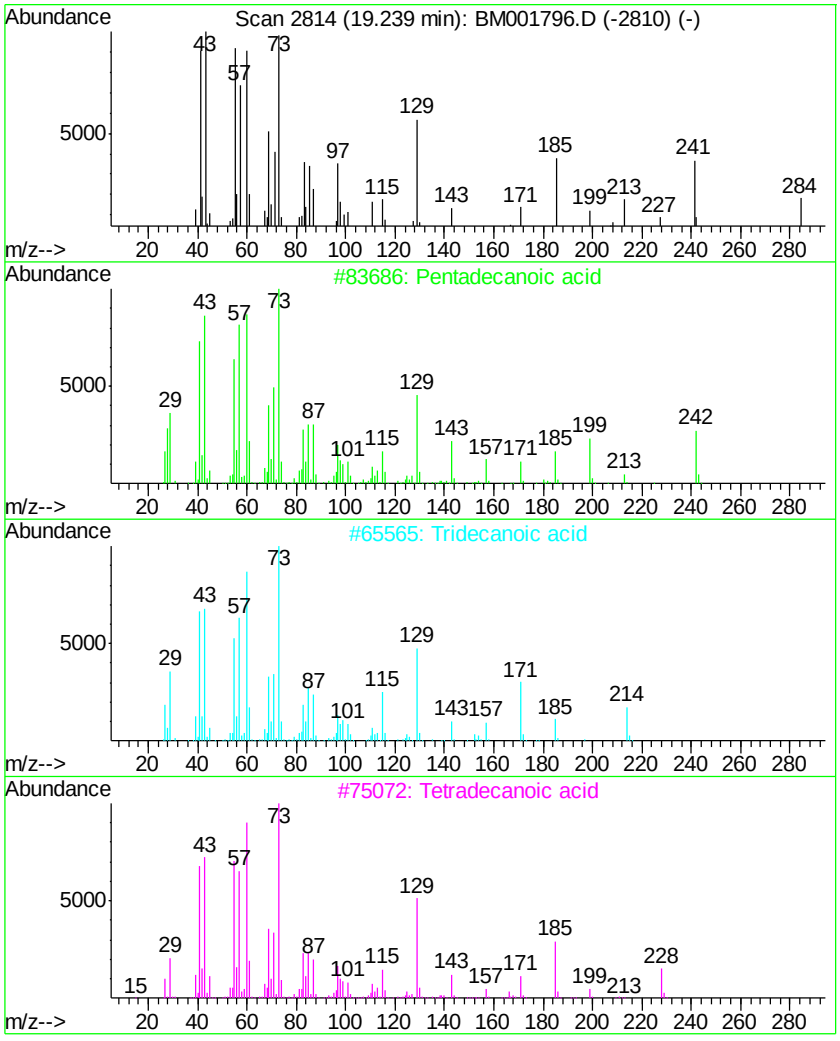
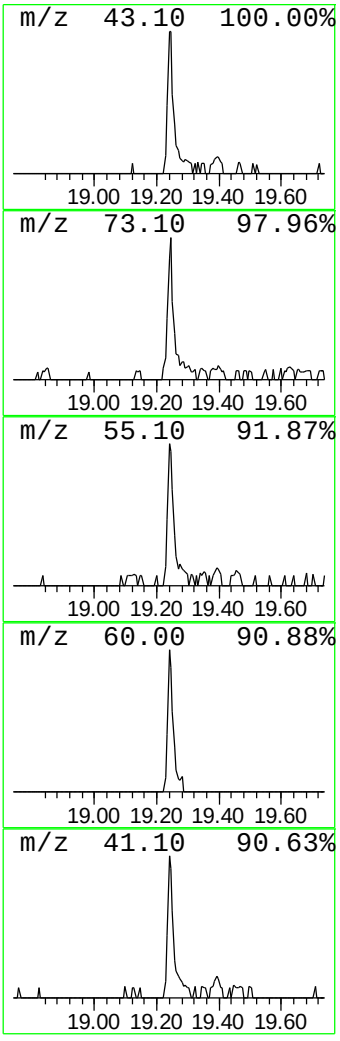
Peak Number 13 Pentadecanoic acid Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.24	2.42 ng/ul	96427	Chrysene-d12	21.26

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentadecanoic acid	242	C15H30O2	001002-84-2	64
2	Tridecanoic acid	214	C13H26O2	000638-53-9	43
3	Tetradecanoic acid	228	C14H28O2	000544-63-8	38
4	n-Decanoic acid	172	C10H20O2	000334-48-5	35
5	n-Decanoic acid	172	C10H20O2	000334-48-5	35





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Acq On : 25 Jun 2015 13:35
Operator : TP/IZ
Sample : G2762-03
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : C01L2

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM061515.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

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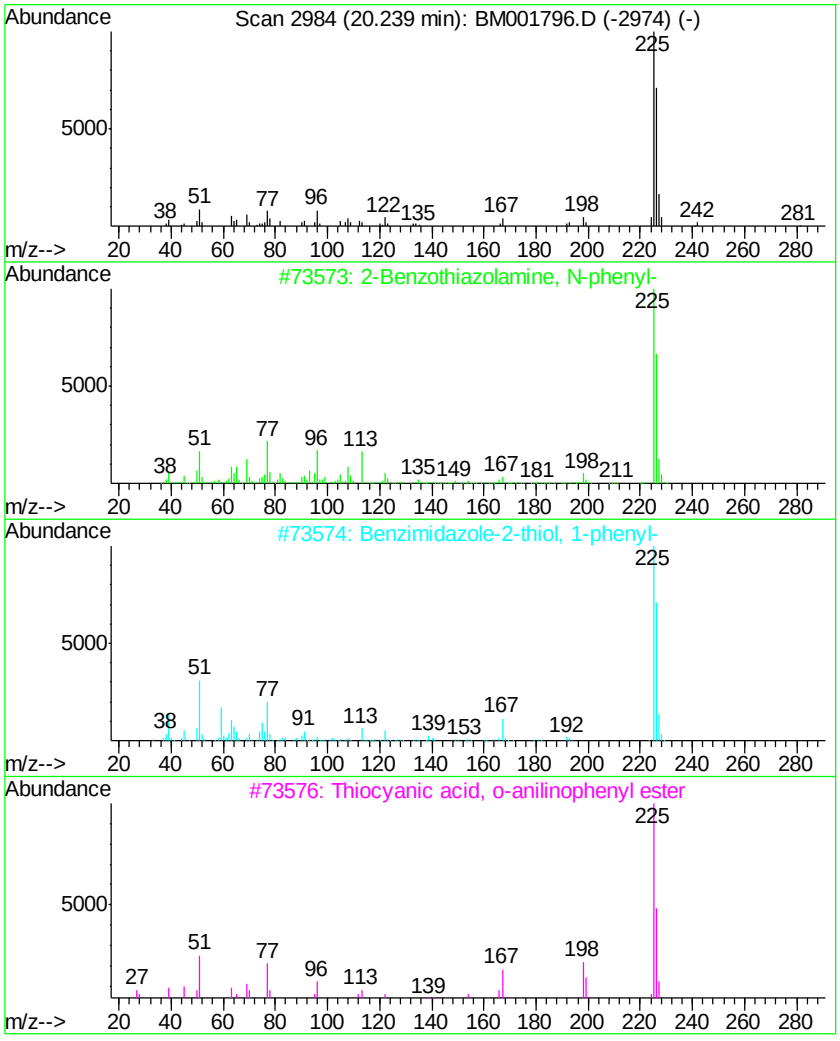
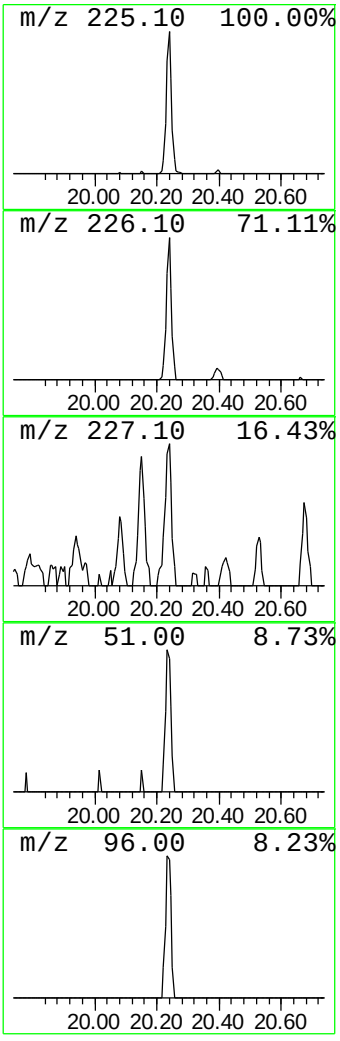
Peak Number 14 2-Benzothiazolamine, N-phenyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.24	2.60 ng/ul	103626	Chrysene-d12	21.26

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Benzothiazolamine, N-phenyl-	226	C13H10N2S	001843-21-6	93
2	Benzimidazole-2-thiol, 1-phenyl-	226	C13H10N2S	004493-32-7	83
3	Thiocyanic acid, o-anilinophenyl...	226	C13H10N2S	015973-81-6	50
4	Drometrizole	225	C13H11N3O	002440-22-4	47
5	2-(p-Fluorophenyl)-1-methylbenzi...	226	C14H11FN2	000724-59-4	46





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Operator : TP/IZ
Sample : G2762-03
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : C01L2

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM061515.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

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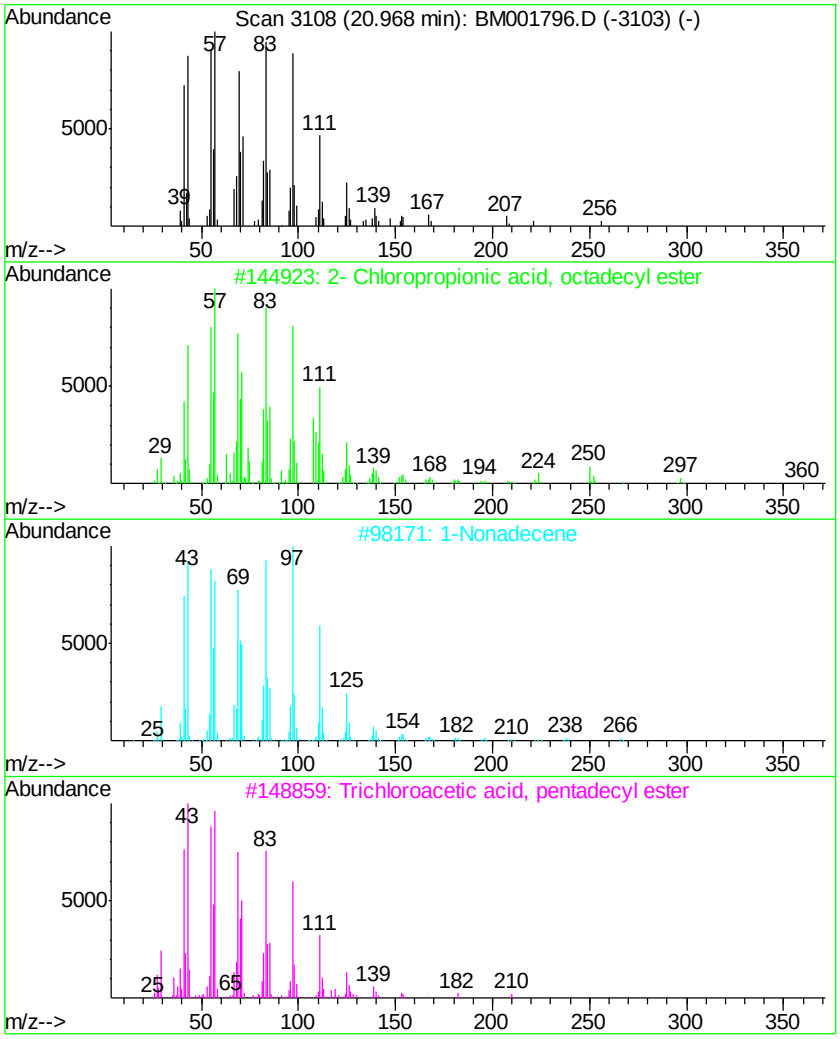
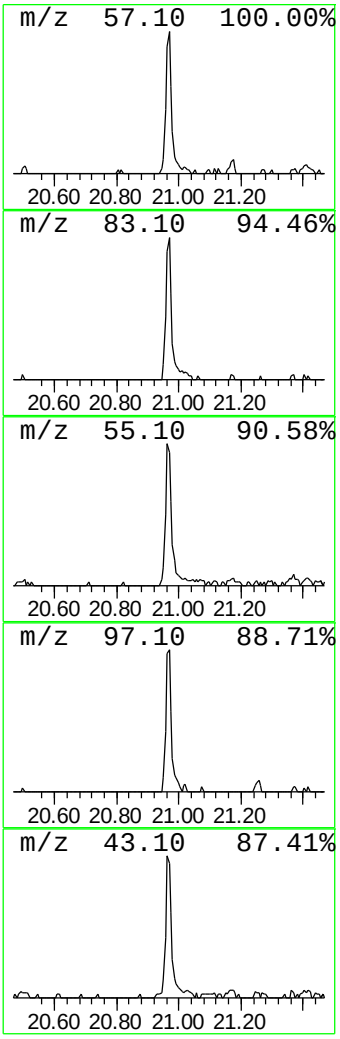
Peak Number 15 2- Chloropropionic acid, oc... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.97	4.65 ng/ul	185151	Chrysene-d12	21.26

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2- Chloropropionic acid, octadec...	360	C21H41ClO2	088104-31-8	94
2	1-Nonadecene	266	C19H38	018435-45-5	91
3	Trichloroacetic acid, pentadecyl...	372	C17H31Cl3O2	074339-53-0	91
4	Trichloroacetic acid, hexadecyl ...	386	C18H33Cl3O2	074339-54-1	91
5	3-Eicosene, (E)-	280	C20H40	074685-33-9	91





Tentatively Identified Compound (LSC) summary

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Data File : BM001796.D
Acq On : 25 Jun 2015 13:35
Operator : TP/IZ
Sample : G2762-03
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : C01L2

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM061515.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

					Internal Standard			
TIC Top Hit name	RT	EstConc	Units	Response	#	RT	Resp	Conc
Butane, 2-methoxy...	2.88	13.4	ng/ul	174360	1	7.66	259782	20.0
Phenol, 3-ethyl-	9.89	10.6	ng/ul	189278	2	10.45	358855	20.0
Phenol, 2,6-dimet...	9.93	8.0	ng/ul	143077	2	10.45	358855	20.0
Phenol, 3,4-dimet...	10.32	7.1	ng/ul	126826	2	10.45	358855	20.0
Benzothiazole	11.10	2.0	ng/ul	36065	2	10.45	358855	20.0
Hydrazine, 1-ethy...	11.27	3.9	ng/ul	69884	2	10.45	358855	20.0
2(3H)-Benzothiazo...	16.07	189.6	ng/ul	5739250	4	17.06	605569	20.0
1,4-Oxathiino[3,2...	16.42	2.5	ng/ul	74647	4	17.06	605569	20.0
n-Hexadecanoic acid	17.98	4.9	ng/ul	148155	4	17.06	605569	20.0
2-Mercaptobenzoth...	18.36	378.6	ng/ul	11462500	4	17.06	605569	20.0
2(3H)-Benzothiazo...	18.50	3.6	ng/ul	109766	4	17.06	605569	20.0
Pentadecanoic acid	19.24	2.4	ng/ul	96427	5	21.26	797056	20.0
2-Benzothiazolami...	20.24	2.6	ng/ul	103626	5	21.26	797056	20.0
2- Chloropropioni...	20.97	4.7	ng/ul	185151	5	21.26	797056	20.0