LSC Area Percent Report

Data Path : Z:\HPCHEM1\BNA G\DATA\BG092116\
Data File : BG024056.D
Acq On : 21 Sep 2016 15:48
Operator : UM/SJ
Sample : H4819-04
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : YORK-SB2-02

Integration Parameters: rteint.p
Integrator: RTE
Smoothing : ON Filtering: 5
Sampling : 1 Min Area: 3 % 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_G\METHODS\8270-BG083116.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.923	9	14	30	rVB	2651178	4034625	100.00%	12.115%
2	5.114	382	387	397	rVB	181530	314238	7.79%	0.944%
3	5.601	464	470	481	rBV	894791	1568830	38.88%	4.711%
4	7.223	739	746	762	rVB	977778	1826133	45.26%	5.484%
5	7.581	800	807	815	rBV	1032606	1788605	44.33%	5.371%
6	8.045	881	886	899	rVB	325960	571953	14.18%	1.717%
7	8.374	935	942	950	rVB	746749	1308191	32.42%	3.928%
8	9.232	1079	1088	1100	rBV	666320	1253445	31.07%	3.764%
9	10.043	1219	1226	1238	rBV5	37056	102035	2.53%	0.306%
10	10.389	1280	1285	1294	rBV4	54548	104966	2.60%	0.315%
11	10.877	1362	1368	1374	rBV	425752	779719	19.33%	2.341%
12	10.930	1374	1377	1383	rVV	105190	166769	4.13%	0.501%
13	11.029	1386	1394	1401	rVB	14851	44846	1.11%	0.135%
14	11.717	1503	1511	1519	rBV2	141525	258362	6.40%	0.776%
15	12.545	1646	1652	1657	rBV	47709	77249	1.91%	0.232%
16	12.762	1683	1689	1700	rBV3	32430	63048	1.56%	0.189%
17	13.332	1778	1786	1794	rBV	1629981	2624150	65.04%	7.880%
18	13.644	1834	1839	1847	rVB	141269	241437	5.98%	0.725%
19	13.896	1875	1882	1887	rBV10	16302	45172	1.12%	0.136%
20	14.037	1901	1906	1911	rBV5	29290	47202	1.17%	0.142%
21	14.166	1923	1928	1934	rVB	334243	482396	11.96%	1.449%
22	14.284	1941	1948	1951	rBV8	22348	48636	1.21%	0.146%
23	14.460	1972	1978	1985	rBV10	25489	57669	1.43%	0.173%
24	14.583	1994	1999	2006	rVB6	32600	47945	1.19%	0.144%
25	14.724	2014	2023	2030	rBV	642397	1045748	25.92%	3.140%
26	14.789	2030	2034	2040	rVB4	25085	42815	1.06%	0.129%
27	15.124	2083	2091	2099	rBV5	27953	62895	1.56%	0.189%
28	15.535	2156	2161	2165	rBV4	73798	121505	3.01%	0.365%
29	15.776	2194	2202	2209	rBV8	26137	67003	1.66%	0.201%
30	16.058	2245	2250	2257	rVB3	60242	105865	2.62%	0.318%
31	16.228	2272	2279	2288	rBV	1201848	1942980	48.16%	5.834%
32	16.410	2306	2310	2312	rBV3	38296	53071	1.32%	0.159%
33	17.203	2442	2445	2448	rBV4	49757	72573	1.80%	0.218%
34	17.309	2459	2463	2469	rVB2	35127	59032	1.46%	0.177%

LSC Area Percent Report

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Operator : UM/SJ
Sample : H4819-04
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : YORK-SB2-02

Integration Parameters: rteint.p
Integrator: RTE
Smoothing : ON Filtering: 5
Sampling : 1 Min Area: 3 % 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_G\METHODS\8270-BG083116.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

35	17.485	2488	2493	2497	rBV	743329	1143285	28.34%	3.433%
36	17.532	2497	2501	2507	rVB	532883	806053	19.98%	2.420%
37	17.626	2513	2517	2521	rVB	83783	117838	2.92%	0.354%
38	18.302	2629	2632	2638	rBV5	76898	138595	3.44%	0.416%
39	18.367	2638	2643	2647	rVV3	311787	524606	13.00%	1.575%
40	18.419	2648	2652	2660	rVB2	134271	288276	7.15%	0.866%
41	18.566	2671	2677	2680	rBV5	108475	242063	6.00%	0.727%
42	18.860	2724	2727	2731	rBV	81899	118086	2.93%	0.355%
43	19.271	2795	2797	2802	rVB4	88891	101137	2.51%	0.304%
44	19.547	2839	2844	2850	rVB	1120352	1527186	37.85%	4.586%
45	19.635	2856	2859	2862	rBV4	133990	185171	4.59%	0.556%
46	19.876	2897	2900	2901	rBV	138726	164299	4.07%	0.493%
47	19.912	2902	2906	2913	rVB	832310	1182070	29.30%	3.550%
48	20.099	2933	2938	2944	rVB	2425188	3266410	80.96%	9.809%
49	21.797	3220	3227	3232	rBV2	831728	2065452	51.19%	6.202%

Sum of corrected areas: 33301635

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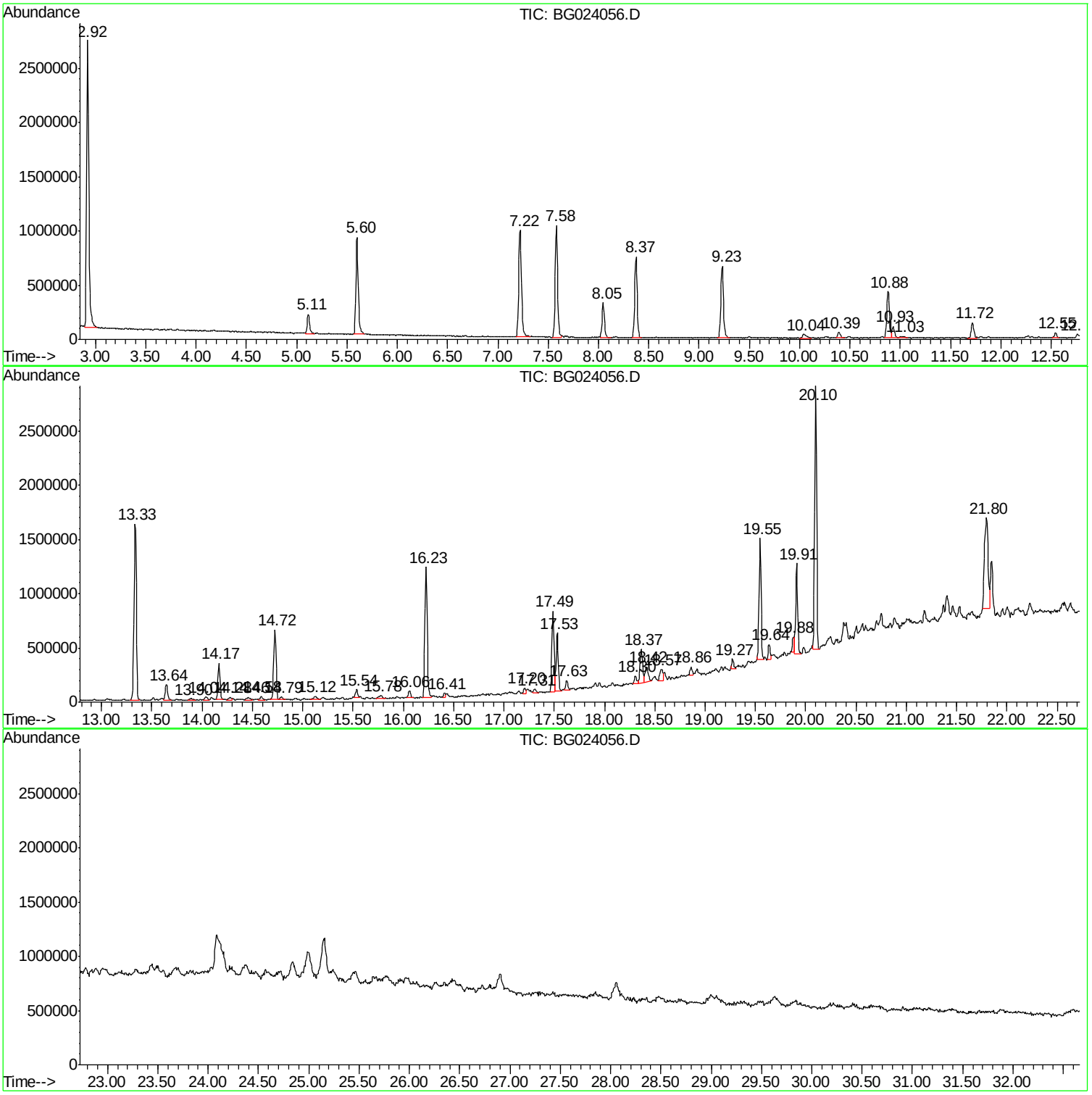
Instrument :
BNA_G
ClientSampleId :
YORK-SB2-02

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG083116.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



Library Search Compound Report

Data Path : Z:\HPCHEM1\BNA G\DATA\BG092116\
Data File : BG024056.D
Acq On : 21 Sep 2016 15:48
Operator : UM/SJ
Sample : H4819-04
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : YORK-SB2-02

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG083116.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

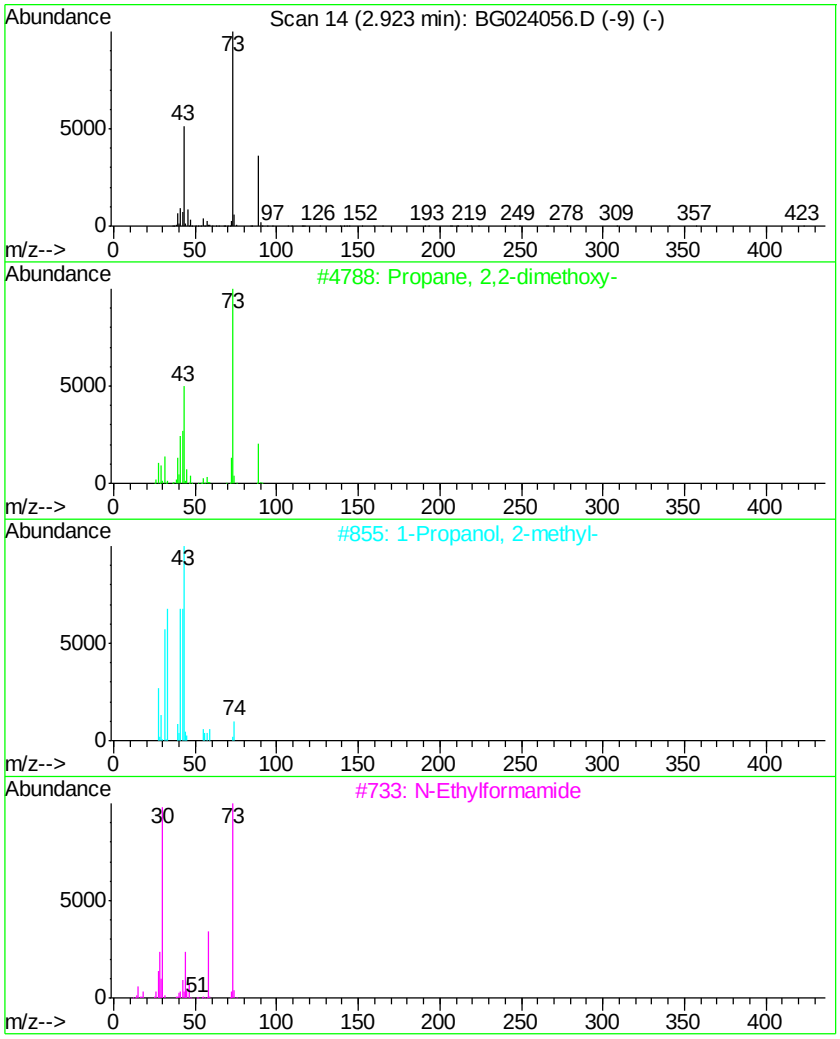
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

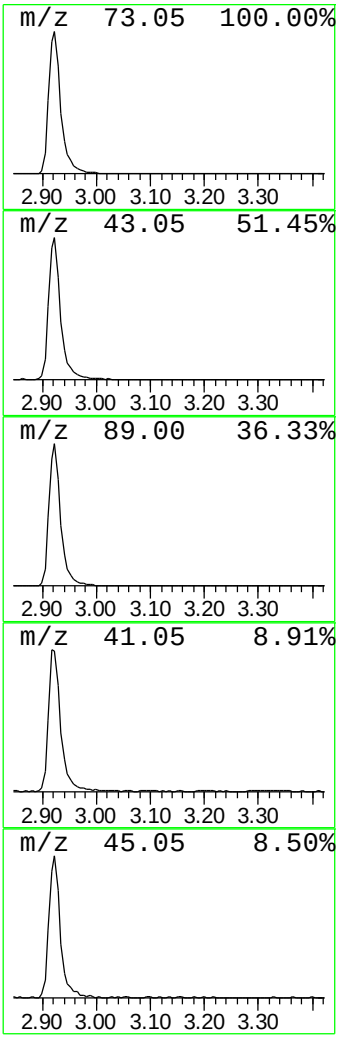
****

Peak Number 1 Propane, 2,2-dimethoxy- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.92	141.08 ng	4034630	1,4-Dichlorobenzene-d4	8.05

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Propane, 2,2-dimethoxy-	104	C5H12O2	000077-76-9	56
2	1-Propanol, 2-methyl-	74	C4H10O	000078-83-1	9
3	N-Ethylformamide	73	C3H7NO	000627-45-2	9
4	1-Hexen-4-ol, 1-chloro-3,5-dimet...	162	C8H15ClO	100244-09-5	9
5	2-Hydroxy-2-methylbutyric acid	118	C5H10O3	003739-30-8	9





Data Path : Z:\HPCHEM1\BNA G\DATA\BG092116\
Data File : BG024056.D
Acq On : 21 Sep 2016 15:48
Operator : UM/SJ
Sample : H4819-04
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : YORK-SB2-02

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG083116.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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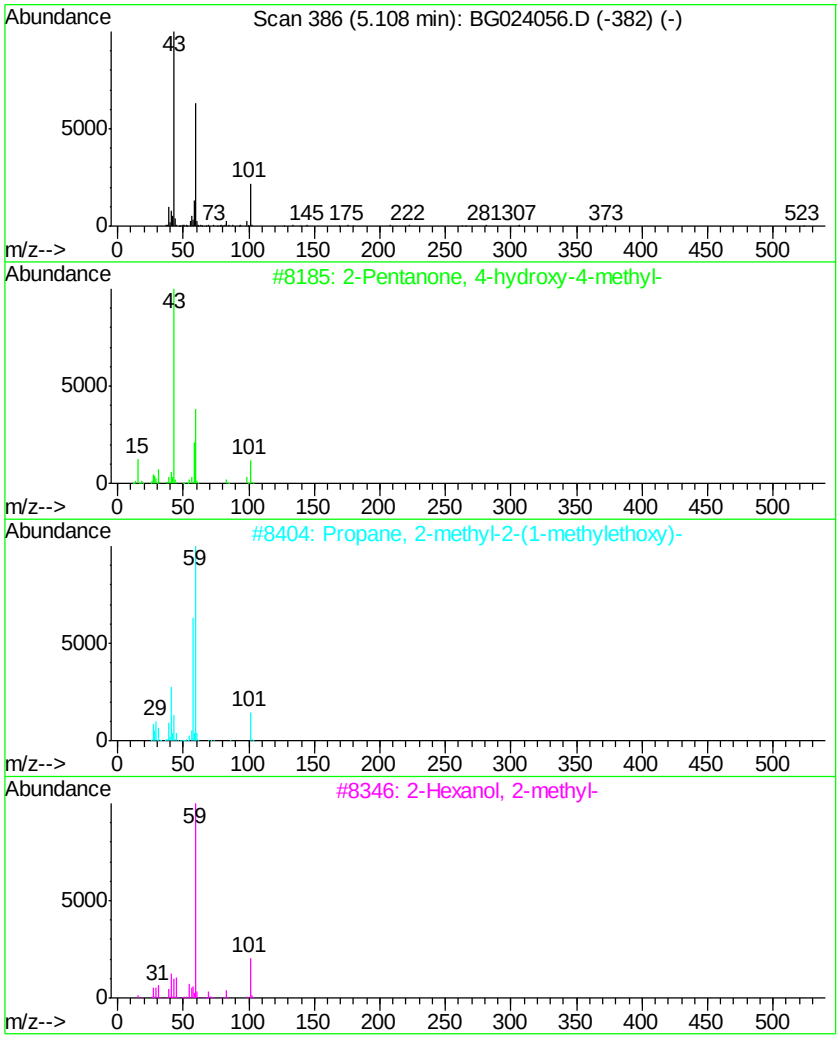
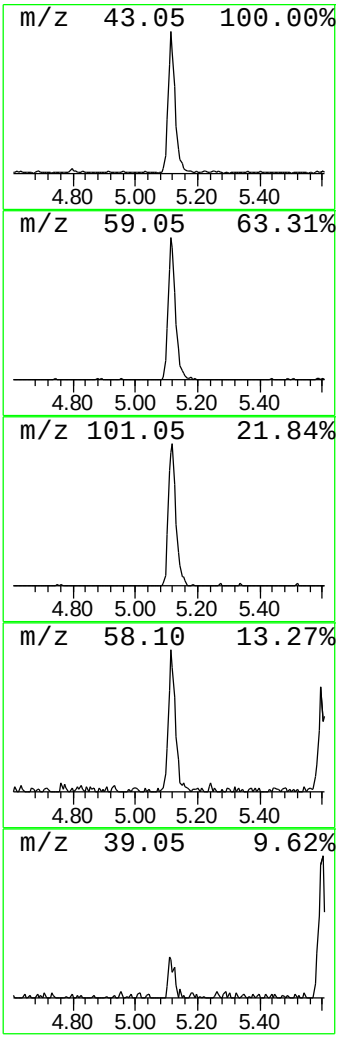
Peak Number 2 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.11	10.99 ng	314238	1,4-Dichlorobenzene-d4	8.05

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	59
2	Propane, 2-methyl-2-(1-methyleth...	116	C7H16O	017348-59-3	50
3	2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	40
4	2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	38
5	Hydrazine, 1,1-bis(1-methylethyl)-	116	C6H16N2	000921-14-2	33





Library Search Compound Report

Data Path : Z:\HPCHEM1\BNA G\DATA\BG092116\
Data File : BG024056.D
Acq On : 21 Sep 2016 15:48
Operator : UM/SJ
Sample : H4819-04
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : YORK-SB2-02

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG083116.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

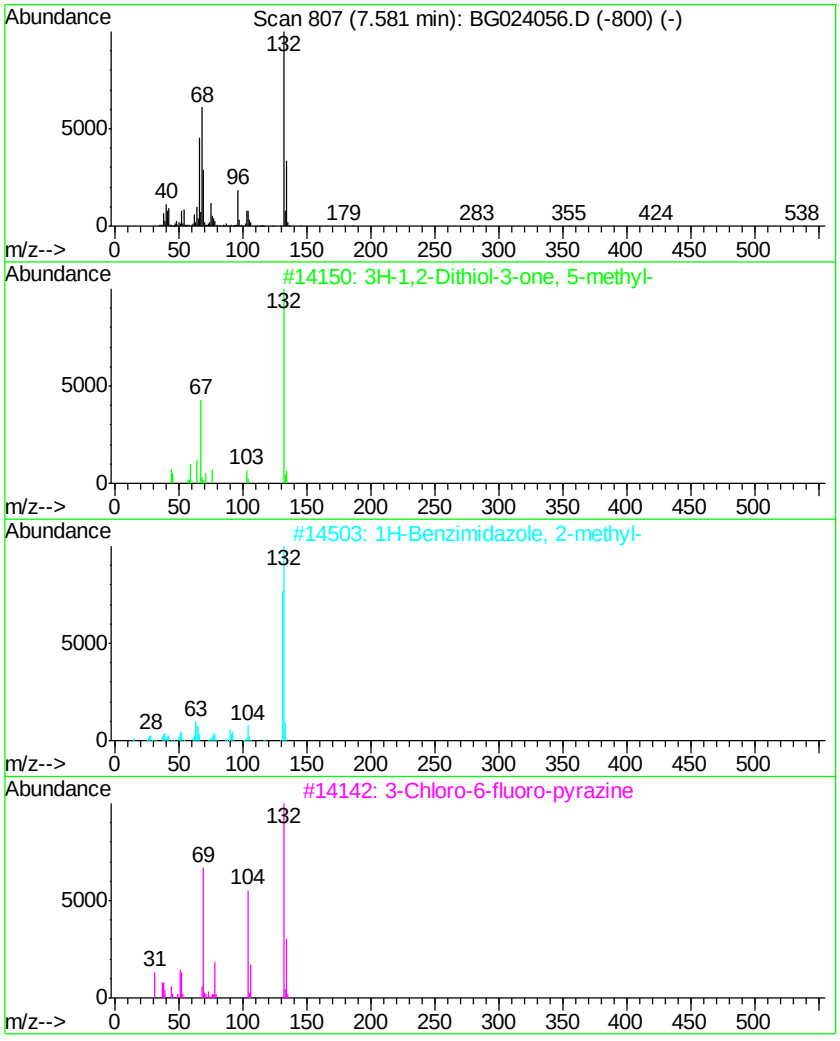
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

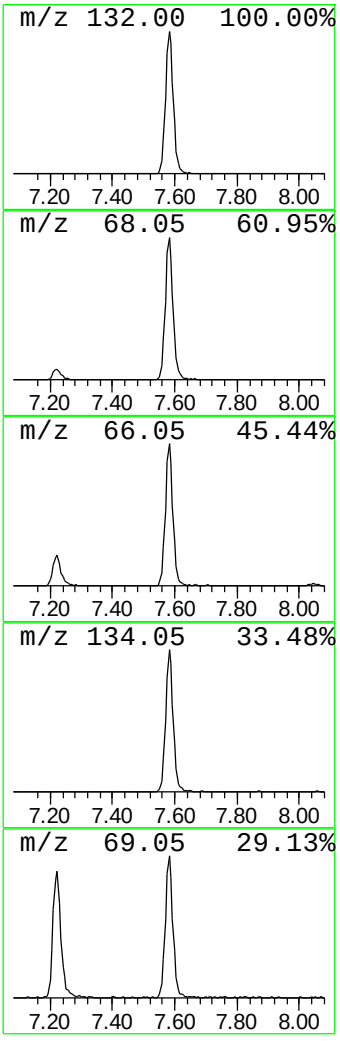
****

Peak Number 3 unknown7.58 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.58	62.54 ng	1788610	1,4-Dichlorobenzene-d4	8.05

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3H-1,2-Dithiol-3-one, 5-methyl-	132	C4H4OS2	003620-08-4	14
2	1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	14
3	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	14
4	1H-Indenol	132	C9H8O	056631-57-3	10
5	Cyclopropanamine, 1-phenyl-	133	C9H11N	1000351-26-2	10





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Operator : UM/SJ
Sample : H4819-04
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : YORK-SB2-02

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG083116.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

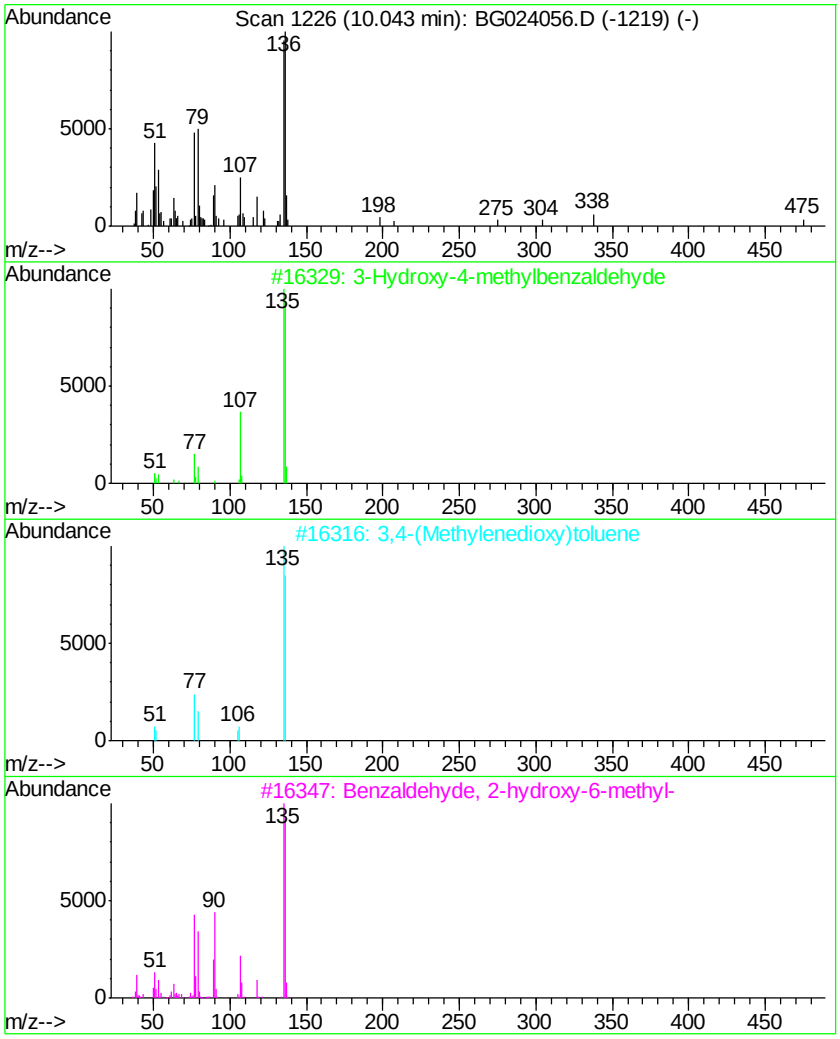
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

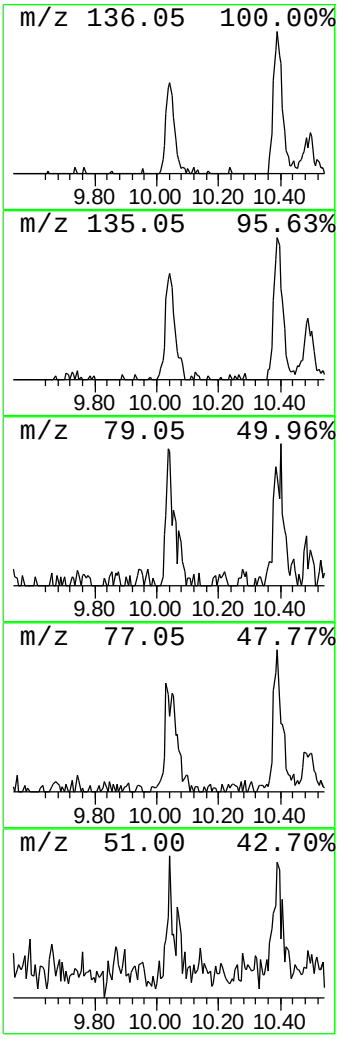
****

Peak Number 5 3-Hydroxy-4-methylbenzaldehyde Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.04	2.62 ng	102035	Naphthalene-d8	10.88

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Hydroxy-4-methylbenzaldehyde	136	C8H8O2	057295-30-4	90
2	3,4-(Methylenedioxy)toluene	136	C8H8O2	007145-99-5	72
3	Benzaldehyde, 2-hydroxy-6-methyl-	136	C8H8O2	018362-36-2	70
4	2-Hydroxy-5-methylbenzaldehyde	136	C8H8O2	000613-84-3	60
5	4-Hydroxy-2-methylbenzaldehyde	136	C8H8O2	041438-18-0	52





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Sample : H4819-04
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : YORK-SB2-02

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG083116.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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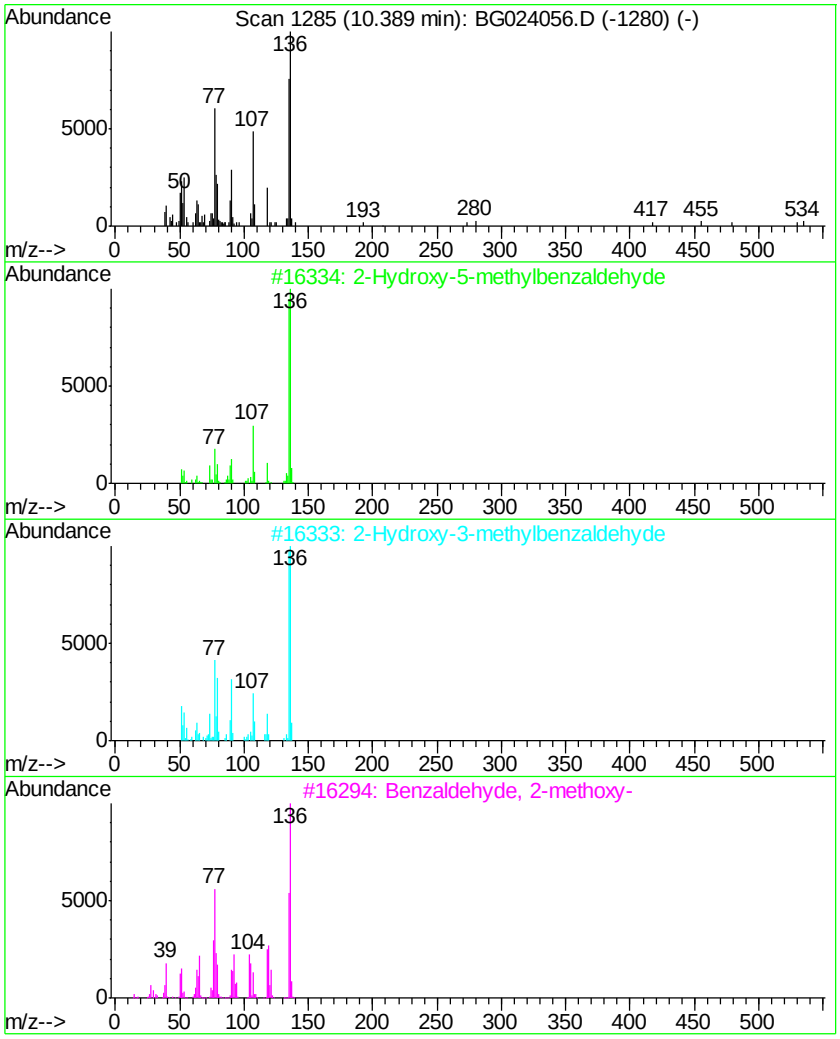
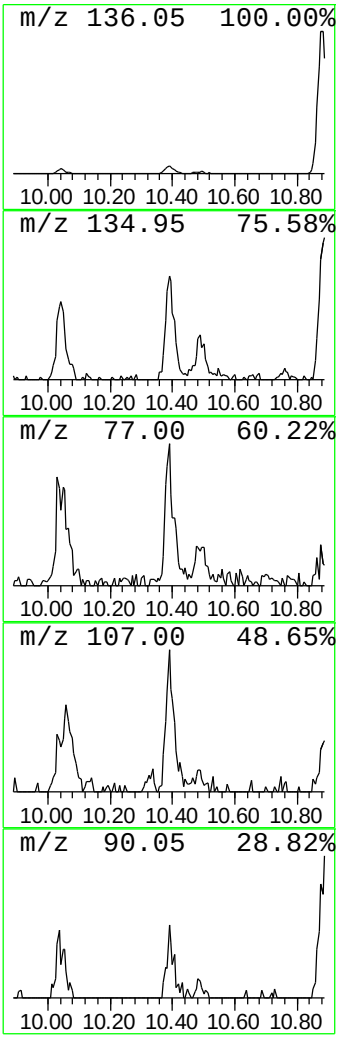
Peak Number 6 2-Hydroxy-5-methylbenzaldehyde Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.39	2.69 ng	104966	Naphthalene-d8	10.88

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Hydroxy-5-methylbenzaldehyde	136	C8H8O2	000613-84-3	72
2	2-Hydroxy-3-methylbenzaldehyde	136	C8H8O2	000824-42-0	68
3	Benzaldehyde, 2-methoxy-	136	C8H8O2	000135-02-4	59
4	3,4-(Methylenedioxy)toluene	136	C8H8O2	007145-99-5	53
5	Benzaldehyde, 2-hydroxy-6-methyl-	136	C8H8O2	018362-36-2	47





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ALS Vial : 7 Sample Multiplier: 1

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ClientSampleId : YORK-SB2-02

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG083116.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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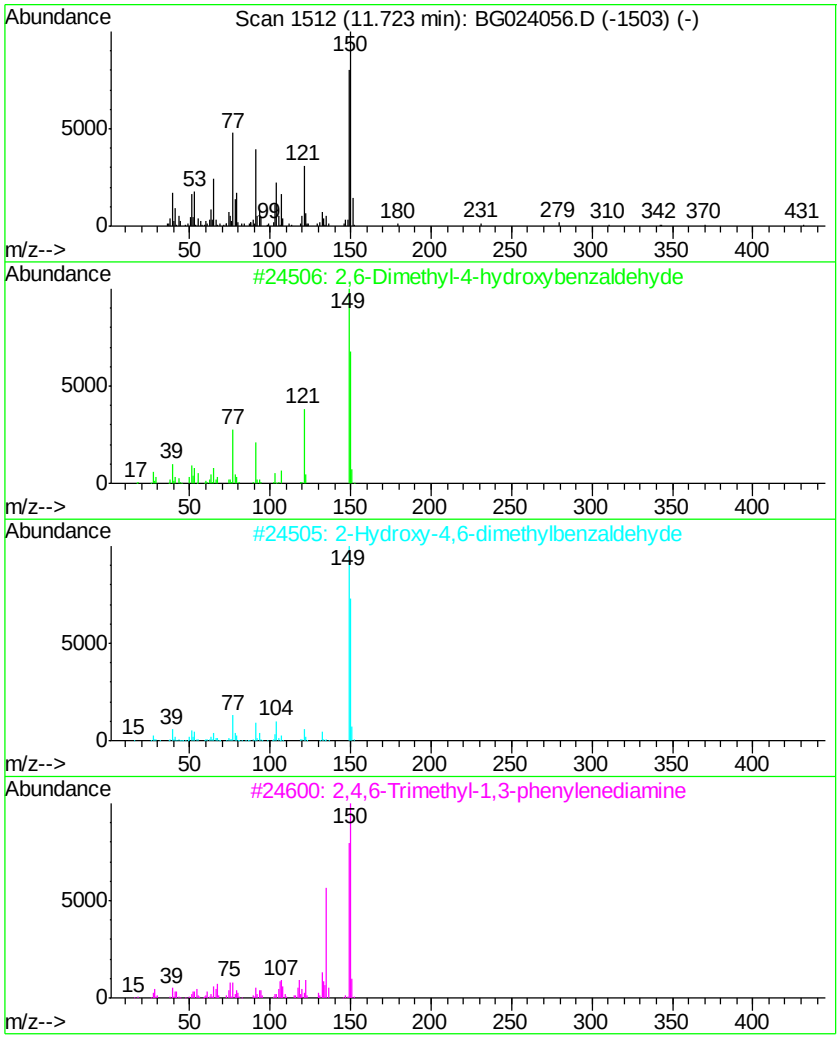
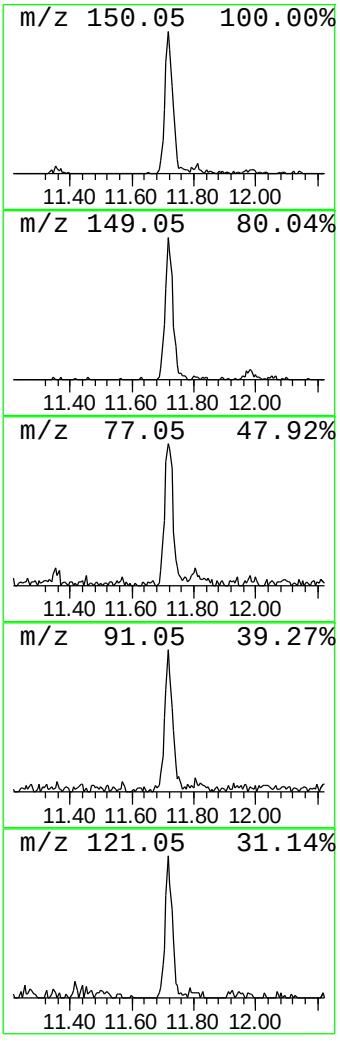
Peak Number 7 2,6-Dimethyl-4-hydroxybenza... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.72	6.63 ng	258362	Naphthalene-d8	10.88

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,6-Dimethyl-4-hydroxybenzaldehyde	150	C9H10O2	070547-87-4	55
2	2-Hydroxy-4,6-dimethylbenzaldehyde	150	C9H10O2	001666-02-0	53
3	2,4,6-Trimethyl-1,3-phenylenedia...	150	C9H14N2	003102-70-3	50
4	4-Pyridinamine, N,N,2,6-tetramet...	150	C9H14N2	129384-12-9	49
5	Benzoic acid, 3,5-dimethyl-	150	C9H10O2	000499-06-9	43





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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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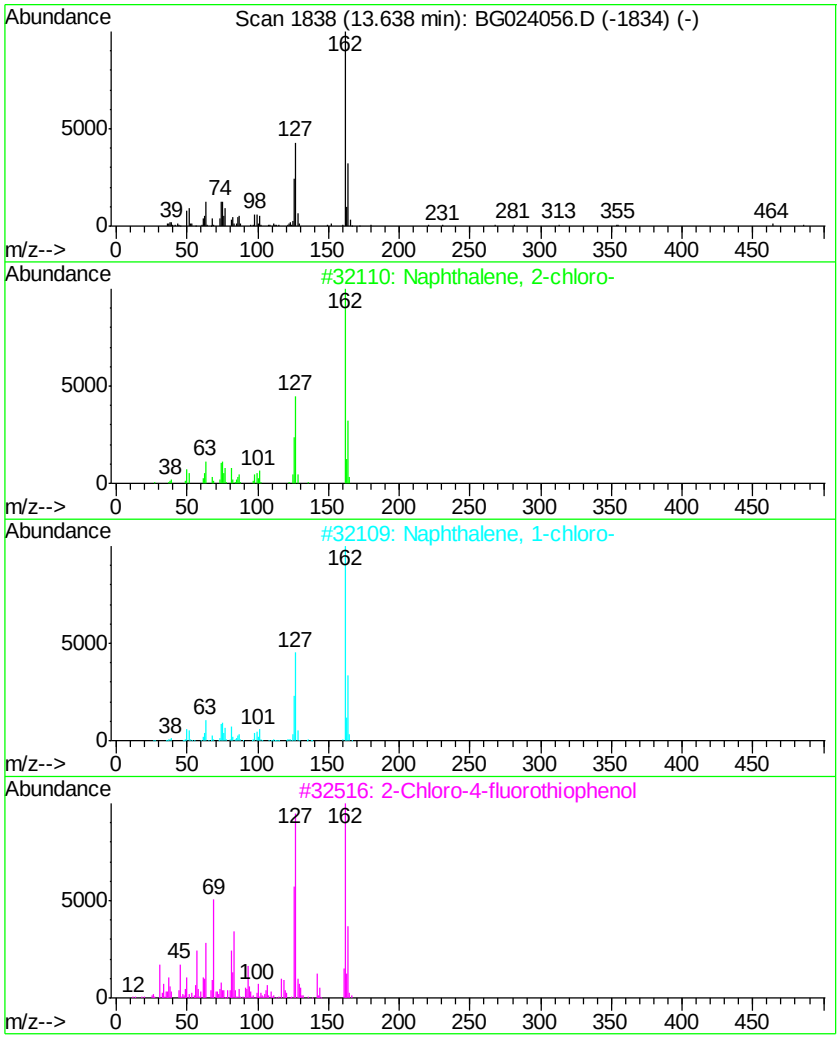
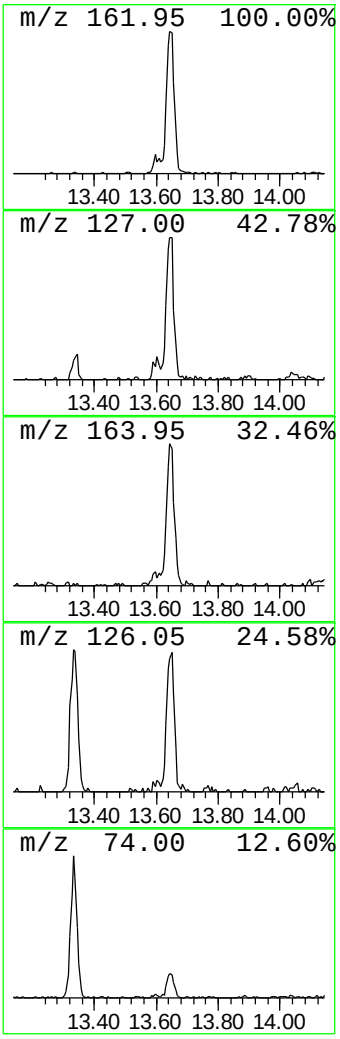
Peak Number 8 Naphthalene, 1-chloro- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.64	4.62 ng	241437	Acenaphthene-d10	14.72

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 2-chloro-	162	C10H7Cl	000091-58-7	97
2	Naphthalene, 1-chloro-	162	C10H7Cl	000090-13-1	96
3	2-Chloro-4-fluorothiophenol	162	C6H4ClFS	175277-99-3	64
4	Pyridin-4-amine, 2,3-dichloro-	162	C5H4Cl2N2	184416-83-9	42
5	6-Nitroindole	162	C8H6N2O2	004769-96-4	38





Data Path : Z:\HPCHEM1\BNA G\DATA\BG092116\
Data File : BG024056.D
Acq On : 21 Sep 2016 15:48
Operator : UM/SJ
Sample : H4819-04
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : YORK-SB2-02

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG083116.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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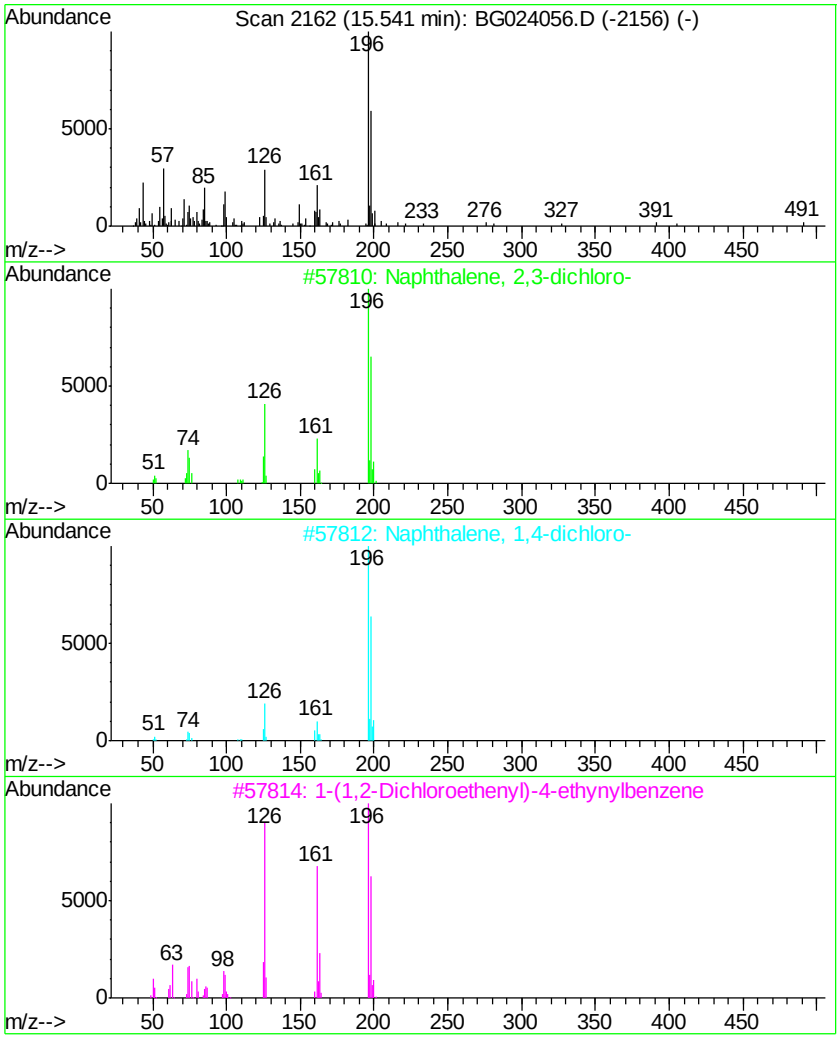
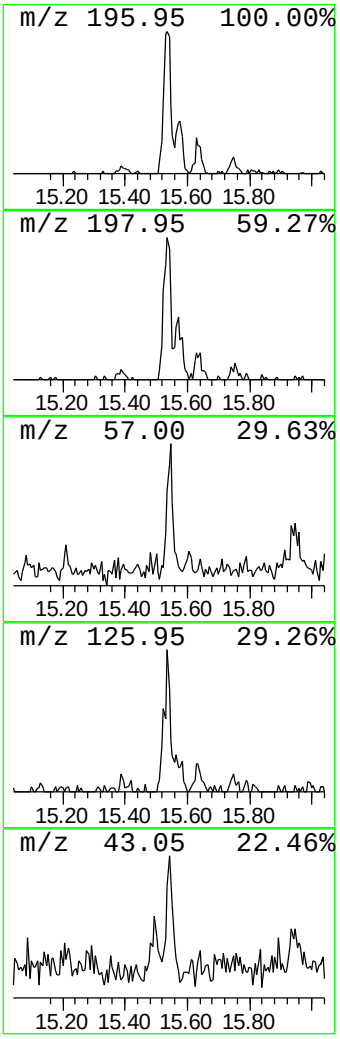
Peak Number 9 Naphthalene, 2,3-dichloro- Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.54	2.32 ng	121505	Acenaphthene-d10	14.72

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 2,3-dichloro-	196	C10H6Cl2	002050-75-1	91
2	Naphthalene, 1,4-dichloro-	196	C10H6Cl2	001825-31-6	91
3	1-(1,2-Dichloroethenyl)-4-ethyny...	196	C10H6Cl2	1000283-76-7	90
4	Naphthalene, 2,7-dichloro-	196	C10H6Cl2	002198-77-8	89
5	Naphthalene, 1,2-dichloro-	196	C10H6Cl2	002050-69-3	87





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Data File : BG024056.D
Acq On : 21 Sep 2016 15:48
Operator : UM/SJ
Sample : H4819-04
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : YORK-SB2-02

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG083116.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

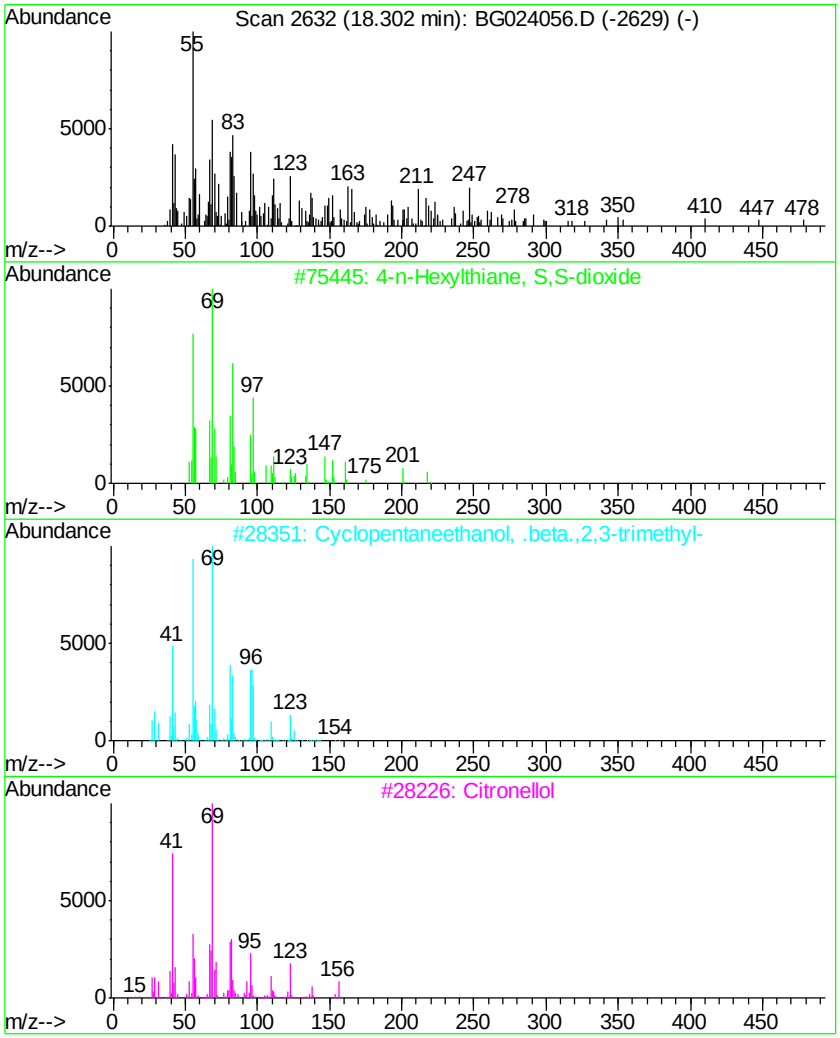
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

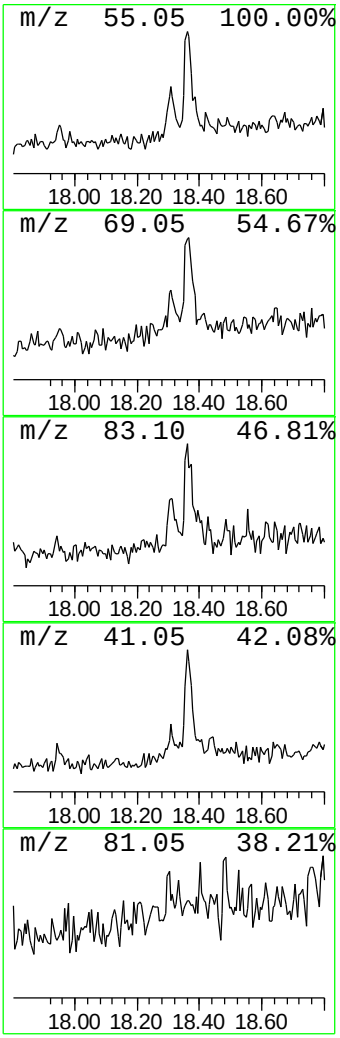
****

Peak Number 11 unknown18.30 Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.30	2.42 ng	138595	Phenanthrene-d10	17.49

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4-n-Hexylthiane, S,S-dioxide	218	C11H22O2S	070928-52-8	15
2	Cyclopentaneethanol, .beta.,2,3-...	156	C10H20O	021721-18-6	12
3	Citronellol	156	C10H20O	000106-22-9	12
4	Palmitoleic acid	254	C16H30O2	000373-49-9	11
5	1,5-Heptadiene, 3-methyl-	110	C8H14	004894-62-6	10





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Operator : UM/SJ
Sample : H4819-04
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : YORK-SB2-02

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG083116.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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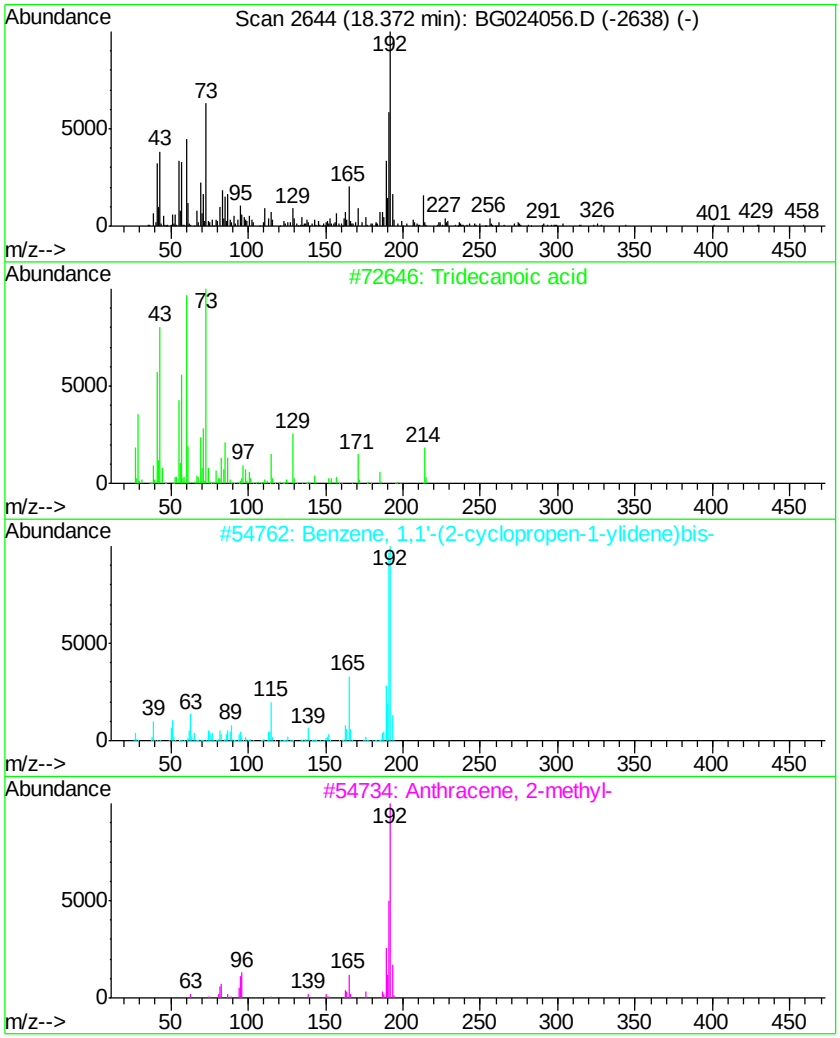
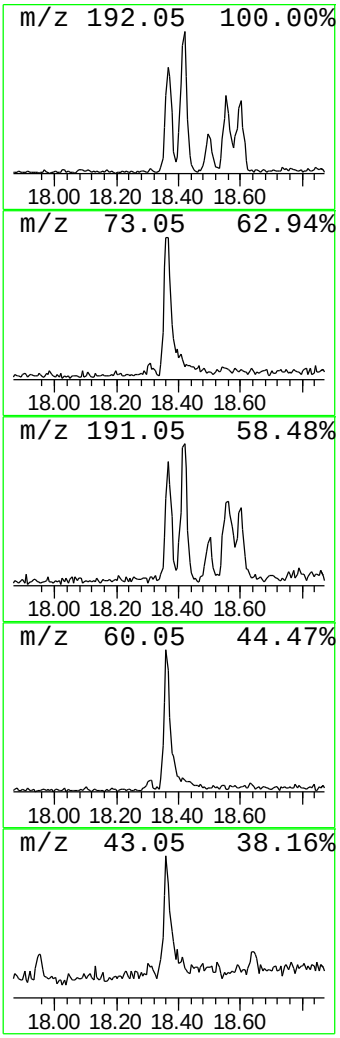
Peak Number 12 Tridecanoic acid Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.37	9.18 ng	524606	Phenanthrene-d10	17.49

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tridecanoic acid	214	C13H26O2	000638-53-9	66
2	Benzene, 1,1'-(2-cyclopropen-1-y...	192	C15H12	022825-21-4	64
3	Anthracene, 2-methyl-	192	C15H12	000613-12-7	60
4	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	51
5	Anthracene, 1-methyl-	192	C15H12	000610-48-0	50





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Operator : UM/SJ
Sample : H4819-04
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : YORK-SB2-02

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG083116.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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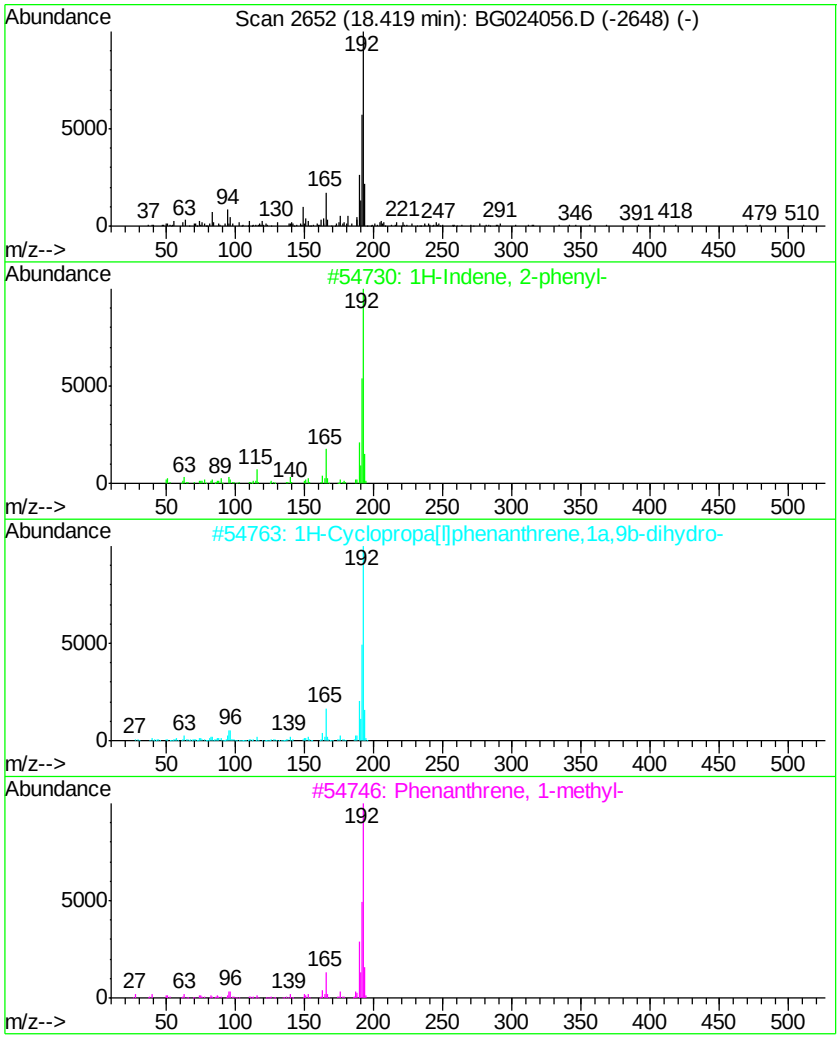
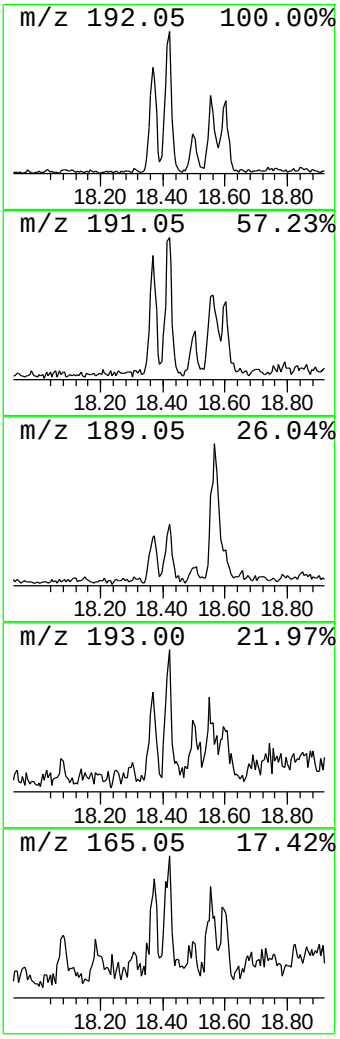
Peak Number 13 1H-Indene, 2-phenyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.42	5.04 ng	288276	Phenanthrene-d10	17.49

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Indene, 2-phenyl-	192	C15H12	004505-48-0	90
2	1H-Cyclopropa[l]phenanthrene,1a,...	192	C15H12	000949-41-7	90
3	Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	90
4	Anthracene, 2-methyl-	192	C15H12	000613-12-7	81
5	Anthracene, 1-methyl-	192	C15H12	000610-48-0	81





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Acq On : 21 Sep 2016 15:48
Operator : UM/SJ
Sample : H4819-04
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : YORK-SB2-02

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG083116.M
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TIC Integration Parameters: LSCINT.P

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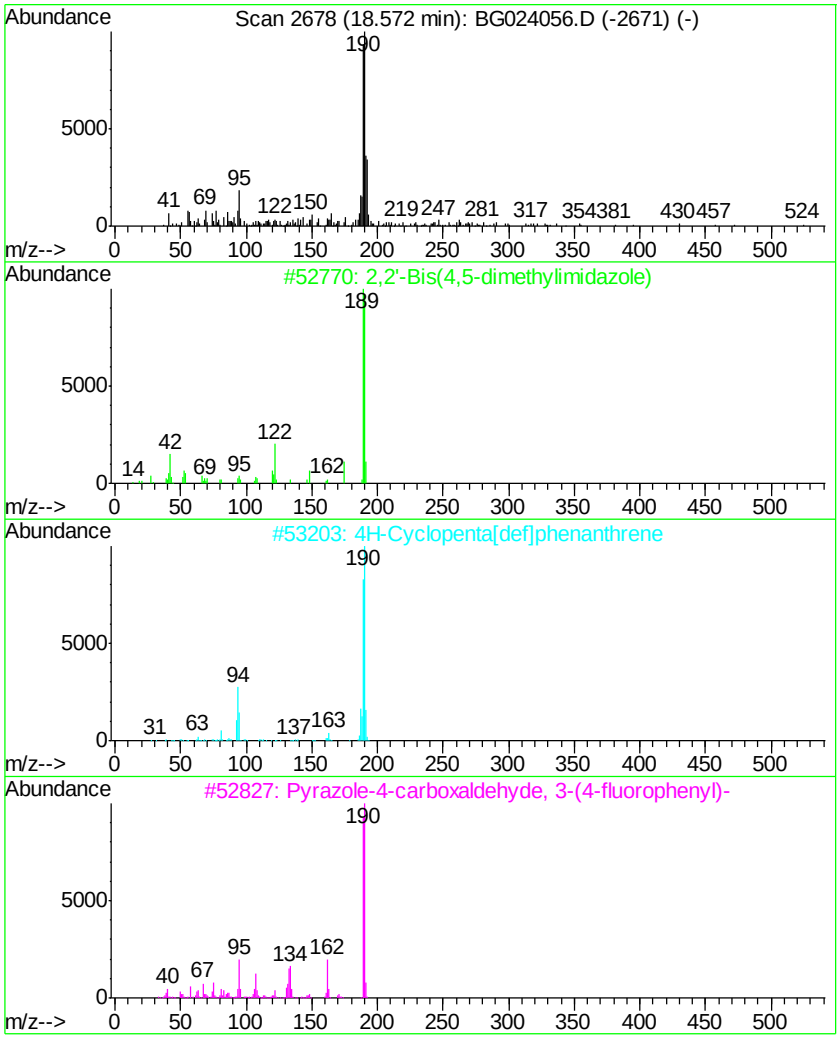
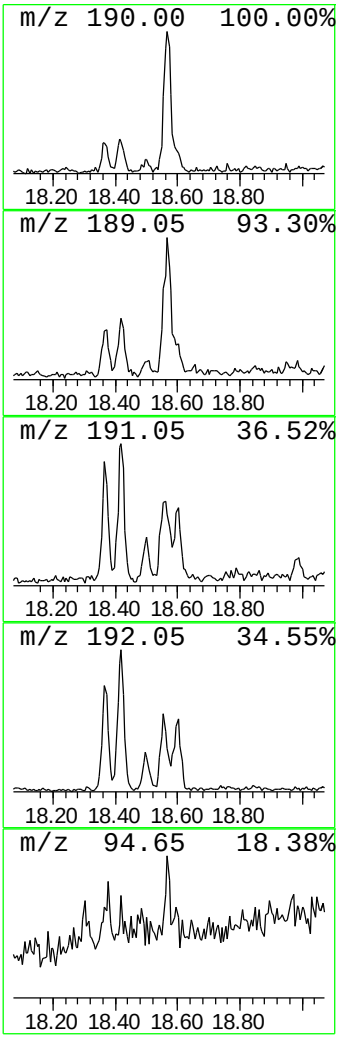
Peak Number 14 unknown18.57 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.57	4.23 ng	242063	Phenanthrene-d10	17.49

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,2'-Bis(4,5-dimethylimidazole)	190	C10H14N4	069286-06-2	49
2	4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	49
3	Pyrazole-4-carboxaldehyde, 3-(4-...	190	C10H7FN2O	306936-57-2	43
4	Phenol, 4-(2-thienylmethyl)-	190	C11H10OS	091680-55-6	38
5	1-Aminocyclopentanecarboxylic ac...	347	C17H30ClNO4	1000329-16-9	27





Library Search Compound Report

Data Path : Z:\HPCHEM1\BNA G\DATA\BG092116\
Data File : BG024056.D
Acq On : 21 Sep 2016 15:48
Operator : UM/SJ
Sample : H4819-04
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : YORK-SB2-02

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG083116.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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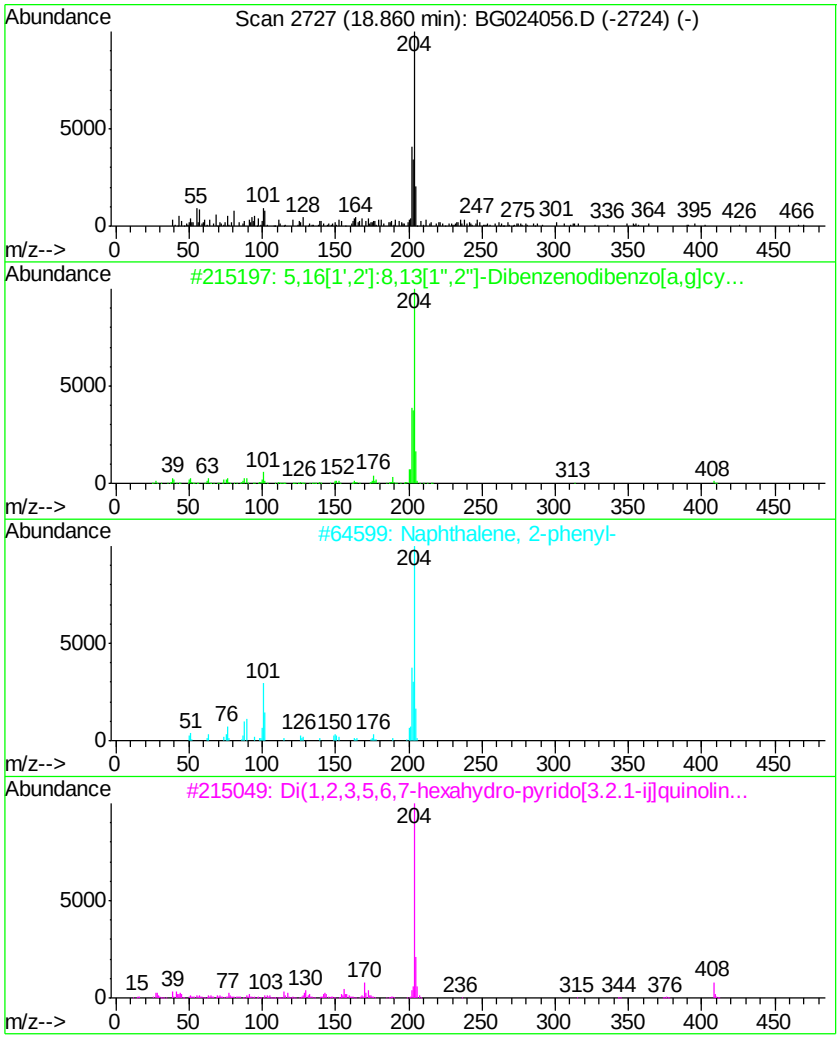
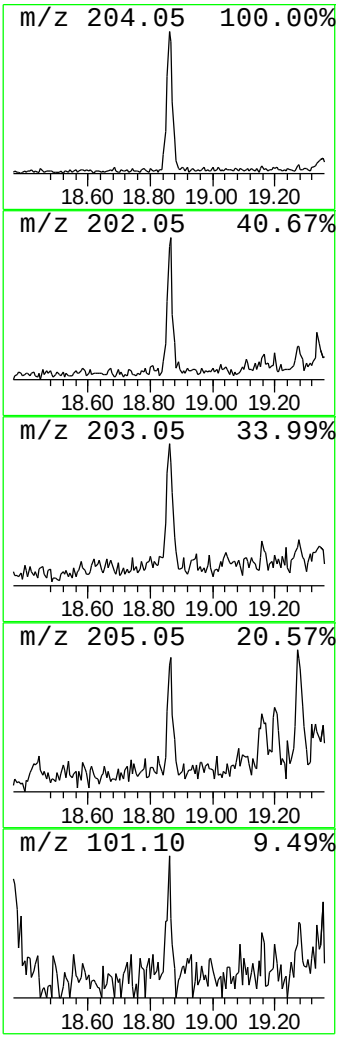
Peak Number 15 5,16[1',2']:8,13[1'',2'']-D... Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.86	2.07 ng	118086	Phenanthrene-d10	17.49

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	5,16[1',2']:8,13[1'',2'']-Dibenz...	408	C32H24	005672-97-9	72
2	Naphthalene, 2-phenyl-	204	C16H12	000612-94-2	52
3	Di(1,2,3,5,6,7-hexahydro-pyrido[...	408	C24H28N2S2	222981-31-9	47
4	Acephenanthrylene, 4,5-dihydro-	204	C16H12	006232-48-0	46
5	Naphthalene, 1-phenyl-	204	C16H12	000605-02-7	38





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ALS Vial : 7 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : YORK-SB2-02

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG083116.M
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TIC Integration Parameters: LSCINT.P

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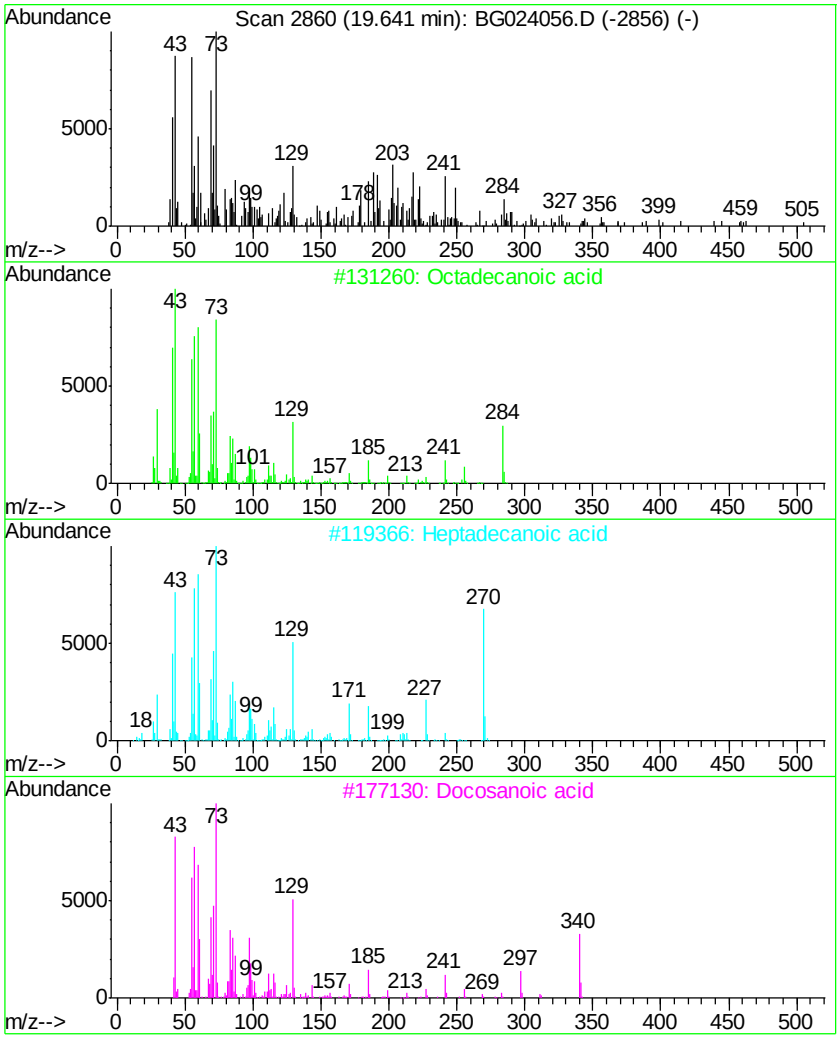
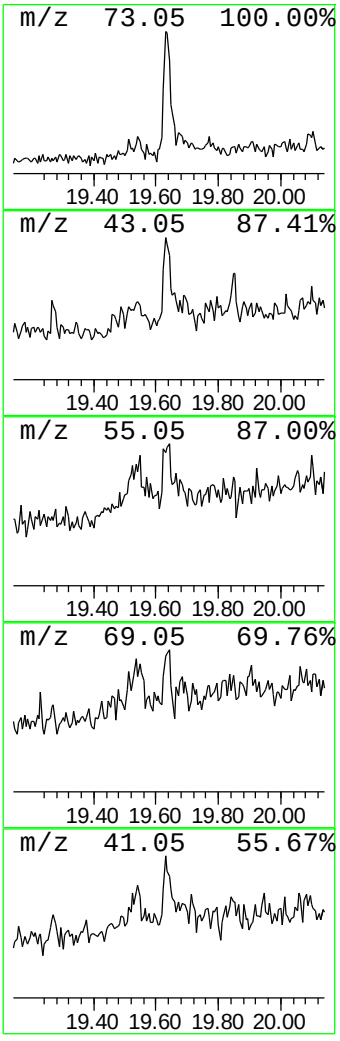
Peak Number 16 Octadecanoic acid Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.64	3.24 ng	185171	Phenanthrene-d10	17.49

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octadecanoic acid	284	C18H36O2	000057-11-4	86
2	Heptadecanoic acid	270	C17H34O2	000506-12-7	42
3	Docosanoic acid	340	C22H44O2	000112-85-6	35
4	11-Bromoundecanoic acid	264	C11H21BrO2	002834-05-1	30
5	Tridecanoic acid	214	C13H26O2	000638-53-9	30





Tentatively Identified Compound (LSC) summary

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Data File : BG024056.D
Acq On : 21 Sep 2016 15:48
Operator : UM/SJ
Sample : H4819-04
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : YORK-SB2-02

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG083116.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT 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
Propane, 2,2-dime...	2.92	141.1	ng	4034630	1	8.05	571953	20.0
2-Pentanone, 4-hy...	5.11	11.0	ng	314238	1	8.05	571953	20.0
unknown7.58	7.58	62.5	ng	1788610	1	8.05	571953	20.0
3-Hydroxy-4-methy...	10.04	2.6	ng	102035	2	10.88	779719	20.0
2-Hydroxy-5-methy...	10.39	2.7	ng	104966	2	10.88	779719	20.0
2,6-Dimethyl-4-hy...	11.72	6.6	ng	258362	2	10.88	779719	20.0
Naphthalene, 1-ch...	13.64	4.6	ng	241437	3	14.72	1045750	20.0
Naphthalene, 2,3-...	15.54	2.3	ng	121505	3	14.72	1045750	20.0
unknown18.30	18.30	2.4	ng	138595	4	17.49	1143290	20.0
Tridecanoic acid	18.37	9.2	ng	524606	4	17.49	1143290	20.0
1H-Indene, 2-phenyl-	18.42	5.0	ng	288276	4	17.49	1143290	20.0
unknown18.57	18.57	4.2	ng	242063	4	17.49	1143290	20.0
5,16[1',2']:8,13[...	18.86	2.1	ng	118086	4	17.49	1143290	20.0
Octadecanoic acid	19.64	3.2	ng	185171	4	17.49	1143290	20.0