LSC Area Percent Report

Data Path : Z:\HPCHEM1\BNA G\DATA\BG092116\
Data File : BG024062.D
Acq On : 21 Sep 2016 19:37
Operator : UM/SJ
Sample : H4821-13 5X
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : YORK-SB1-4

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	2.928	11	15	25	rVB	583697	822405	50.14%	5.472%
2	4.515	282	285	295	rVB4	23147	38326	2.34%	0.255%
3	5.125	384	389	394	rBV	65380	106374	6.49%	0.708%
4	5.613	467	472	482	rBV	94803	168028	10.24%	1.118%
5	7.017	706	711	713	rBV5	14919	19511	1.19%	0.130%
6	7.234	742	748	761	rBV2	128914	268188	16.35%	1.784%
7	7.593	803	809	815	rBV2	123559	217559	13.26%	1.448%
8	8.051	880	887	896	rBV2	371283	658726	40.16%	4.383%
9	8.380	936	943	950	rBV2	121041	227279	13.86%	1.512%
10	9.238	1083	1089	1100	rBV2	96775	197501	12.04%	1.314%
11	10.888	1363	1370	1377	rBV	496992	895333	54.58%	5.957%
12	13.338	1777	1787	1793	rBV	256795	419411	25.57%	2.791%
13	13.579	1822	1828	1835	rBV7	21543	38981	2.38%	0.259%
14	14.031	1899	1905	1912	rBV2	246759	376169	22.93%	2.503%
15	14.178	1924	1930	1935	rBV2	77927	128497	7.83%	0.855%
16	14.342	1952	1958	1965	rBV	476195	654520	39.90%	4.355%
17	14.425	1967	1972	1975	rBV	51317	78508	4.79%	0.522%
18	14.495	1980	1984	1987	rVV2	112044	154222	9.40%	1.026%
19	14.519	1987	1988	1994	rVB4	37413	52194	3.18%	0.347%
20	14.642	2004	2009	2012	rBV2	37682	52773	3.22%	0.351%
21	14.683	2012	2016	2019	rVV3	93175	140467	8.56%	0.935%
22	14.730	2019	2024	2030	rVV	747145	1203050	73.34%	8.004%
23	14.777	2030	2032	2042	rVV9	27022	71615	4.37%	0.476%
24	14.865	2042	2047	2051	rVB3	32020	46409	2.83%	0.309%
25	14.959	2059	2063	2066	rBV4	35371	56555	3.45%	0.376%
26	14.995	2066	2069	2074	rVB3	81123	111850	6.82%	0.744%
27	15.388	2128	2136	2142	rBV2	170844	247616	15.10%	1.648%
28	15.688	2180	2187	2190	rBV6	27773	45764	2.79%	0.304%
29	15.729	2190	2194	2200	rVB5	59043	88056	5.37%	0.586%
30	15.846	2210	2214	2218	rVB7	14766	21525	1.31%	0.143%
31	16.158	2262	2267	2269	rBV5	23589	34092	2.08%	0.227%
32	16.216	2273	2277	2288	rVB6	78468	178480	10.88%	1.188%
33	16.457	2316	2318	2323	rVB6	20489	24163	1.47%	0.161%
34	16.598	2334	2342	2346	rBV9	14686	39811	2.43%	0.265%

LSC Area Percent Report

Data Path : Z:\HPCHEM1\BNA G\DATA\BG092116\
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Acq On : 21 Sep 2016 19:37
Operator : UM/SJ
Sample : H4821-13 5X
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : YORK-SB1-4

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	16.839	2378	2383	2388	rVB5	163219	254169	15.50%	1.691%
36	16.945	2398	2401	2406	rVB5	52847	64337	3.92%	0.428%
37	17.004	2406	2411	2414	rVV7	20512	35080	2.14%	0.233%
38	17.033	2414	2416	2420	rVB4	17816	22521	1.37%	0.150%
39	17.491	2488	2494	2502	rBV2	826015	1291912	78.76%	8.596%
40	18.225	2613	2619	2624	rBV	1209474	1640276	100.00%	10.914%
41	18.308	2631	2633	2639	rBV6	32039	44839	2.73%	0.298%
42	18.366	2639	2643	2647	rBV7	52112	86208	5.26%	0.574%
43	20.099	2934	2938	2945	rBV	406073	548557	33.44%	3.650%
44	21.497	3173	3176	3188	rVB	272026	399329	24.35%	2.657%
45	21.803	3222	3228	3238	rVB	892612	1495080	91.15%	9.947%
46	25.157	3793	3799	3811	rVB2	446568	1263467	77.03%	8.406%

Sum of corrected areas: 15029733

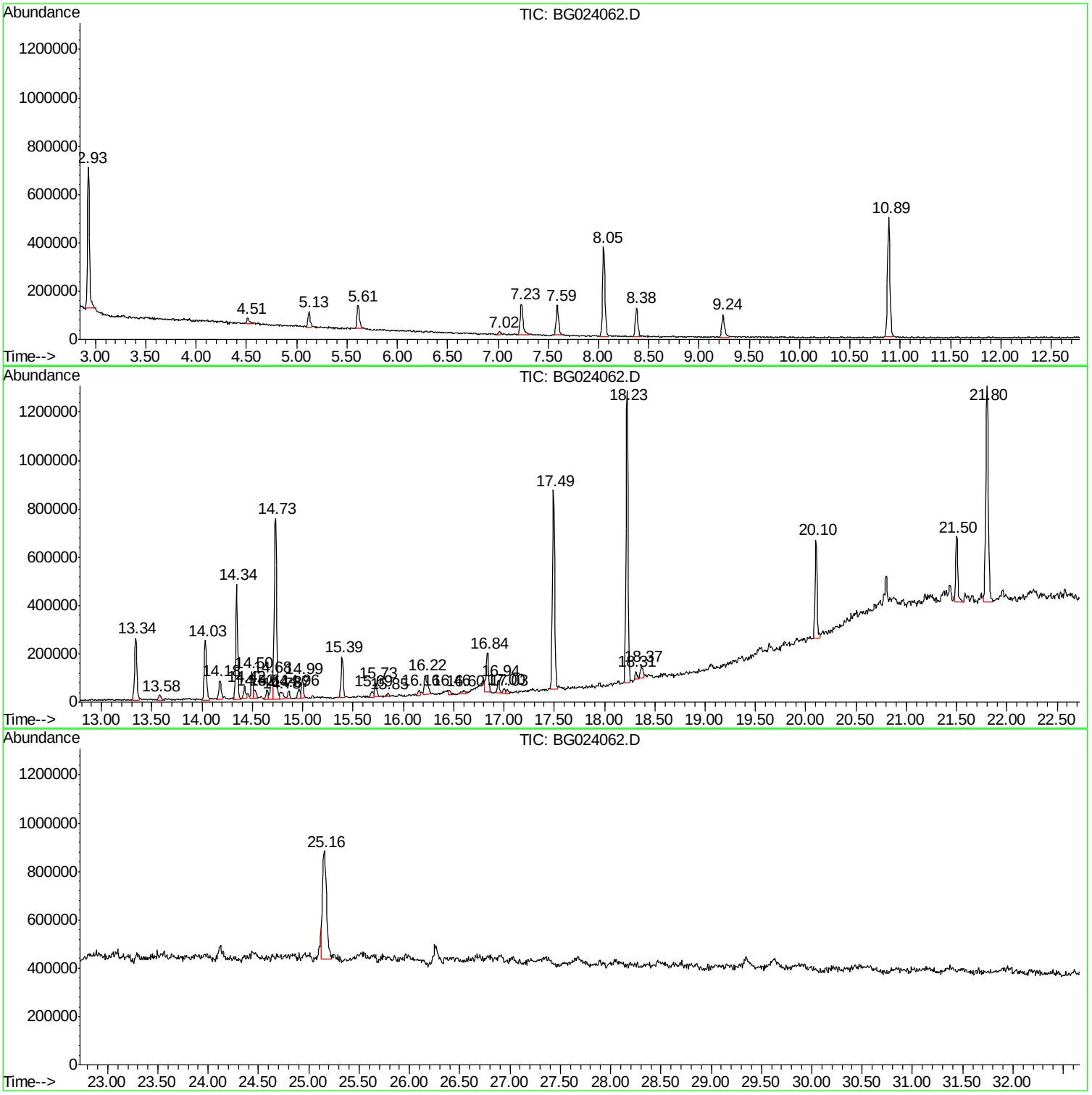
LSC Report - Integrated Chromatogram

Data Path : Z:\HPCHEM1\BNA G\DATA\BG092116\
Data File : BG024062.D
Acq On : 21 Sep 2016 19:37
Operator : UM/SJ
Sample : H4821-13 5X
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
YORK-SB1-4

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



Data Path : Z:\HPCHEM1\BNA G\DATA\BG092116\
Data File : BG024062.D
Acq On : 21 Sep 2016 19:37
Operator : UM/SJ
Sample : H4821-13 5X
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : YORK-SB1-4

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

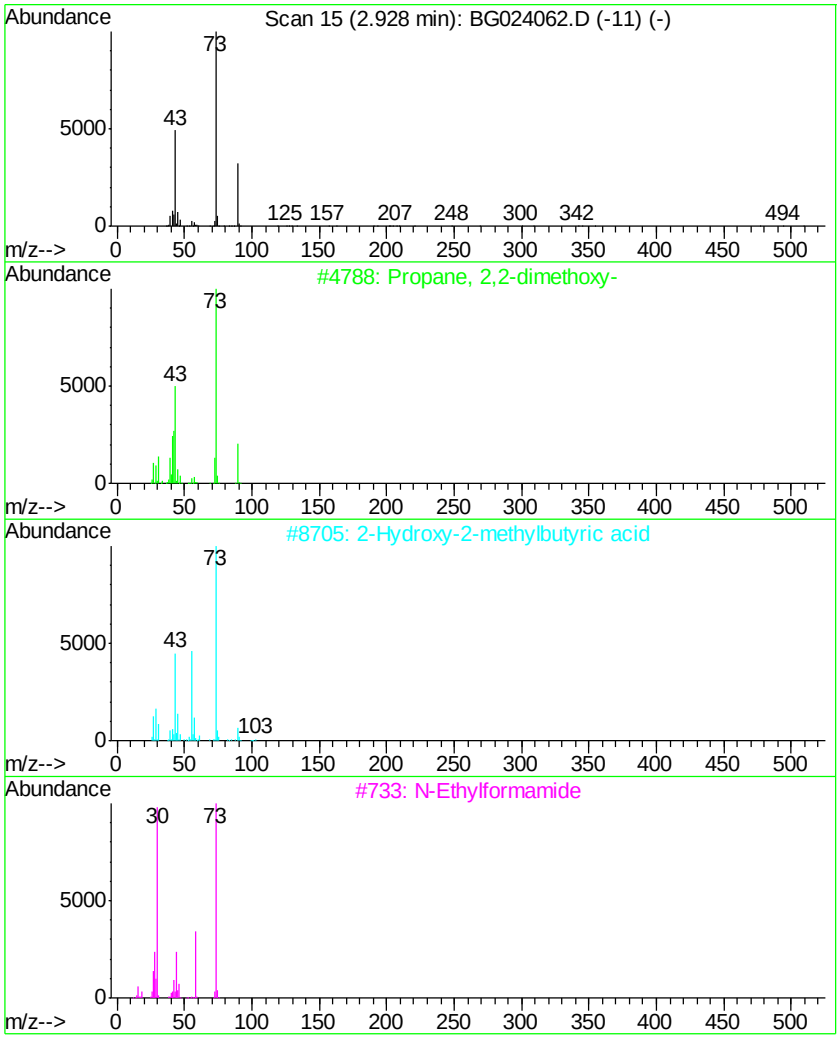
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

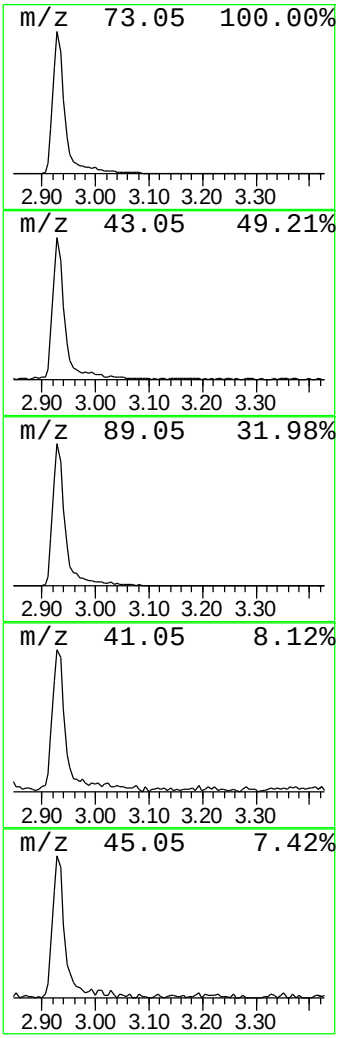
****

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.93	24.97 ng	822405	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	50
2	2-Hydroxy-2-methylbutyric acid	118	C5H10O3	003739-30-8	33
3	N-Ethylformamide	73	C3H7NO	000627-45-2	9
4	Pentane, 3-methoxy-	102	C6H14O	036839-67-5	9
5	Methane, isothiocyanato-	73	C2H3NS	000556-61-6	7





Data Path : Z:\HPCHEM1\BNA G\DATA\BG092116\
Data File : BG024062.D
Acq On : 21 Sep 2016 19:37
Operator : UM/SJ
Sample : H4821-13 5X
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : YORK-SB1-4

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

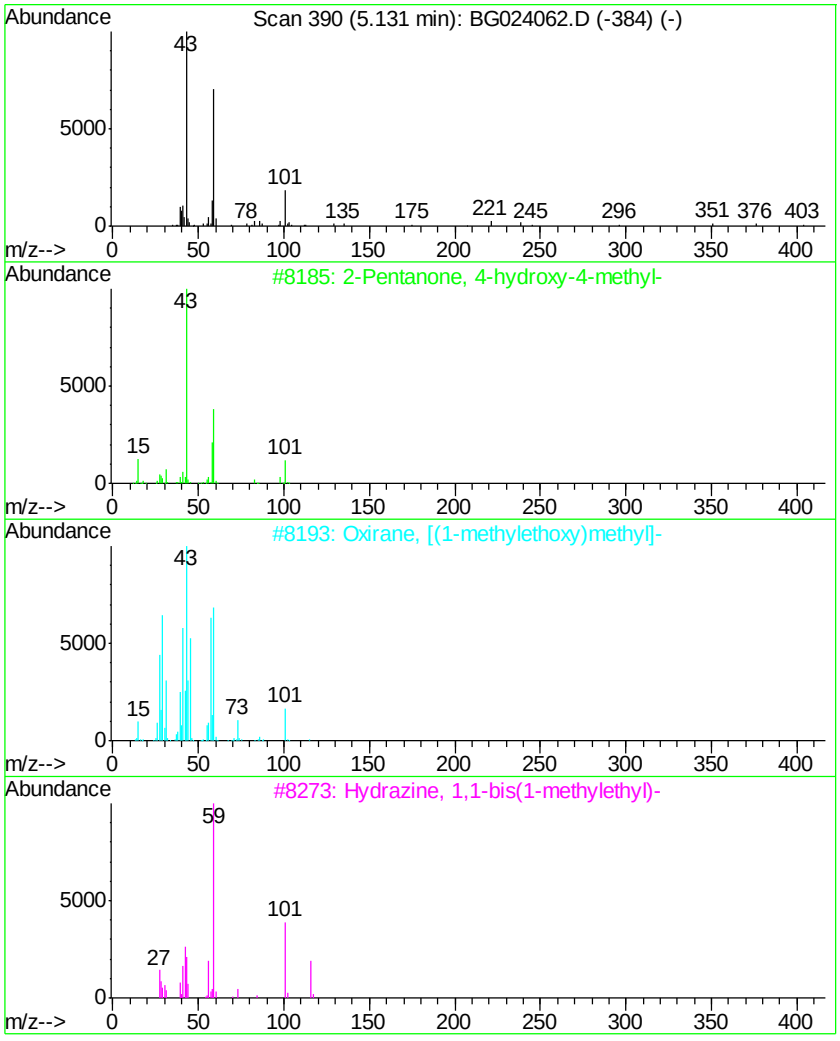
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

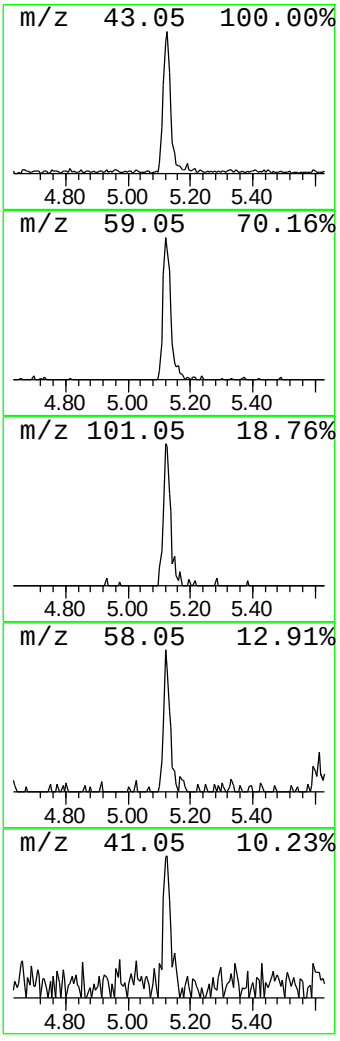
****

Peak Number 2 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.13	3.23 ng	106374	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	42
2	Oxirane, [(1-methylethoxy)methyl]-	116	C6H12O2	004016-14-2	40
3	Hydrazine, 1,1-bis(1-methylethyl)-	116	C6H16N2	000921-14-2	28
4	3-Octanol	130	C8H18O	000589-98-0	28
5	Butanedioic acid, diazo-, dimeth...	172	C6H8N2O4	055514-36-8	28





Data Path : Z:\HPCHEM1\BNA G\DATA\BG092116\
Data File : BG024062.D
Acq On : 21 Sep 2016 19:37
Operator : UM/SJ
Sample : H4821-13 5X
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : YORK-SB1-4

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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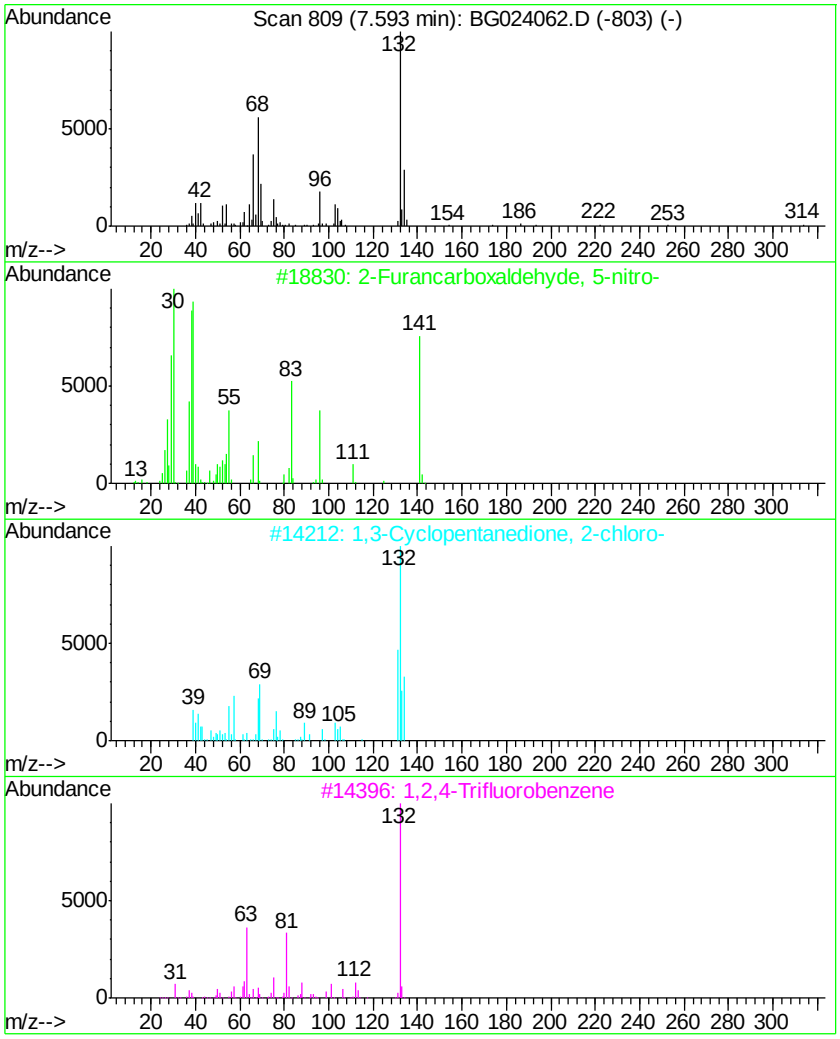
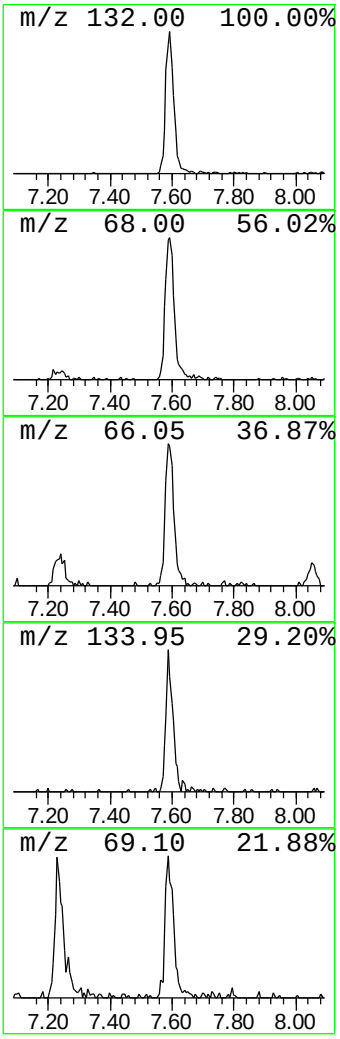
Peak Number 3 unknown7.59 Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.59	6.61 ng	217559	1,4-Dichlorobenzene-d4	8.05

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Furancarboxaldehyde, 5-nitro-	141	C5H3NO4	000698-63-5	35
2	1,3-Cyclopentanedione, 2-chloro-	132	C5H5ClO2	014203-19-1	32
3	1,2,4-Trifluorobenzene	132	C6H3F3	000367-23-7	14
4	1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	12
5	3H-1,2-Dithiol-3-one, 5-methyl-	132	C4H4OS2	003620-08-4	12





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Misc :
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Instrument : BNA_G
ClientSampleId : YORK-SB1-4

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG083116.M
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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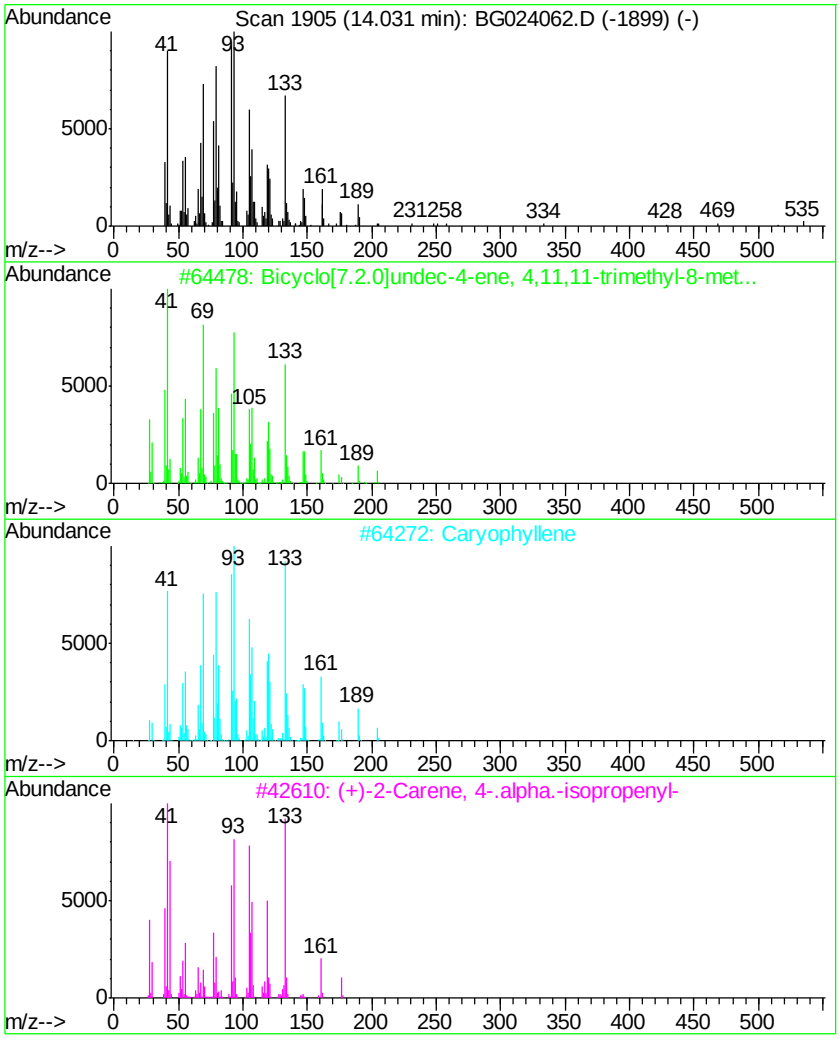
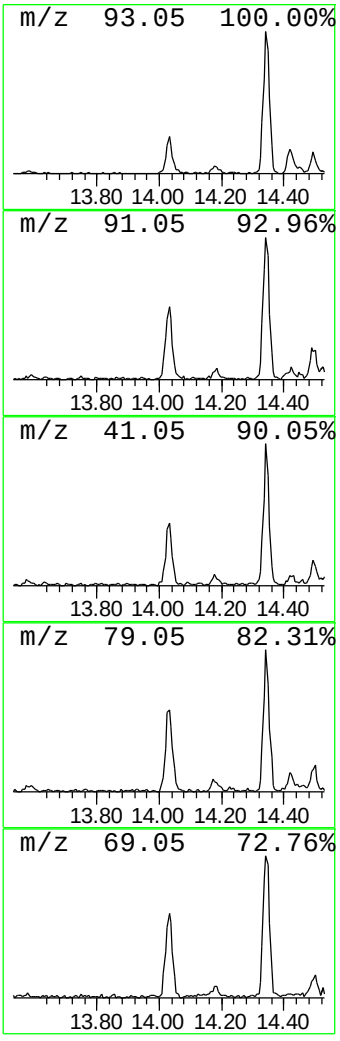
Peak Number 5 Bicyclo[7.2.0]undec-4-ene, ... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.03	6.25 ng	376169	Acenaphthene-d10	14.73

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Bicyclo[7.2.0]undec-4-ene, 4,11,...	204	C15H24	000118-65-0	95
2	Caryophyllene	204	C15H24	000087-44-5	95
3	(+)-2-Carene, 4-.alpha.-isoprope...	176	C13H20	1000151-26-0	87
4	Isocaryophillene	204	C15H24	1000140-07-2	72
5	Aromandendrene	204	C15H24	000489-39-4	62





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Operator : UM/SJ
Sample : H4821-13 5X
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : YORK-SB1-4

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

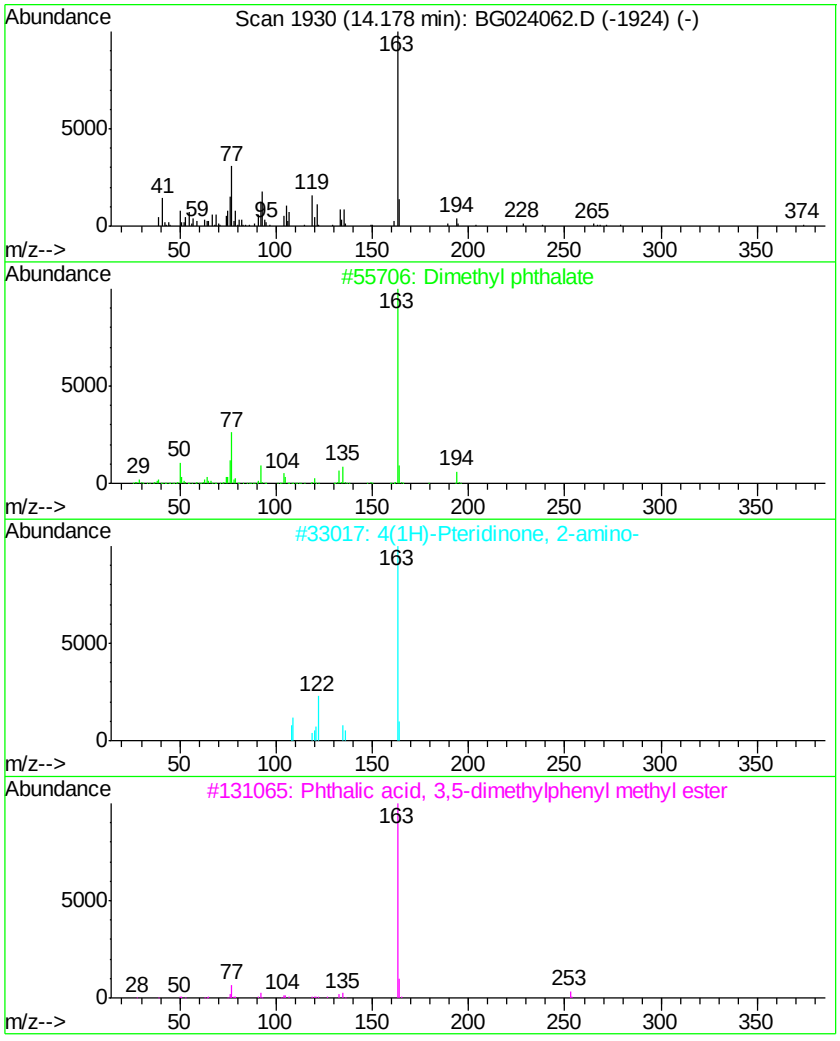
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

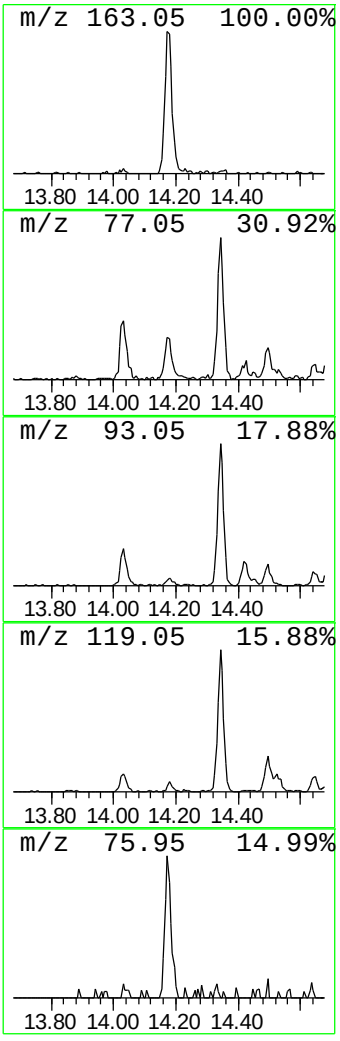
****

Peak Number 6 Dimethyl phthalate Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.18	2.14 ng	128497	Acenaphthene-d10	14.73

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Dimethyl phthalate	194	C10H10O4	000131-11-3	87
2	4(1H)-Pteridinone, 2-amino-	163	C6H5N5O	002236-60-4	78
3	Phthalic acid, 3,5-dimethylphenyl...	284	C17H16O4	1000315-70-8	72
4	Hexyl methyl phthalate	264	C15H20O4	1000373-89-7	64
5	Benzoic acid, 2-(1-oxopropyl)-, ...	192	C11H12O3	032025-37-9	64





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Sample : H4821-13 5X
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : YORK-SB1-4

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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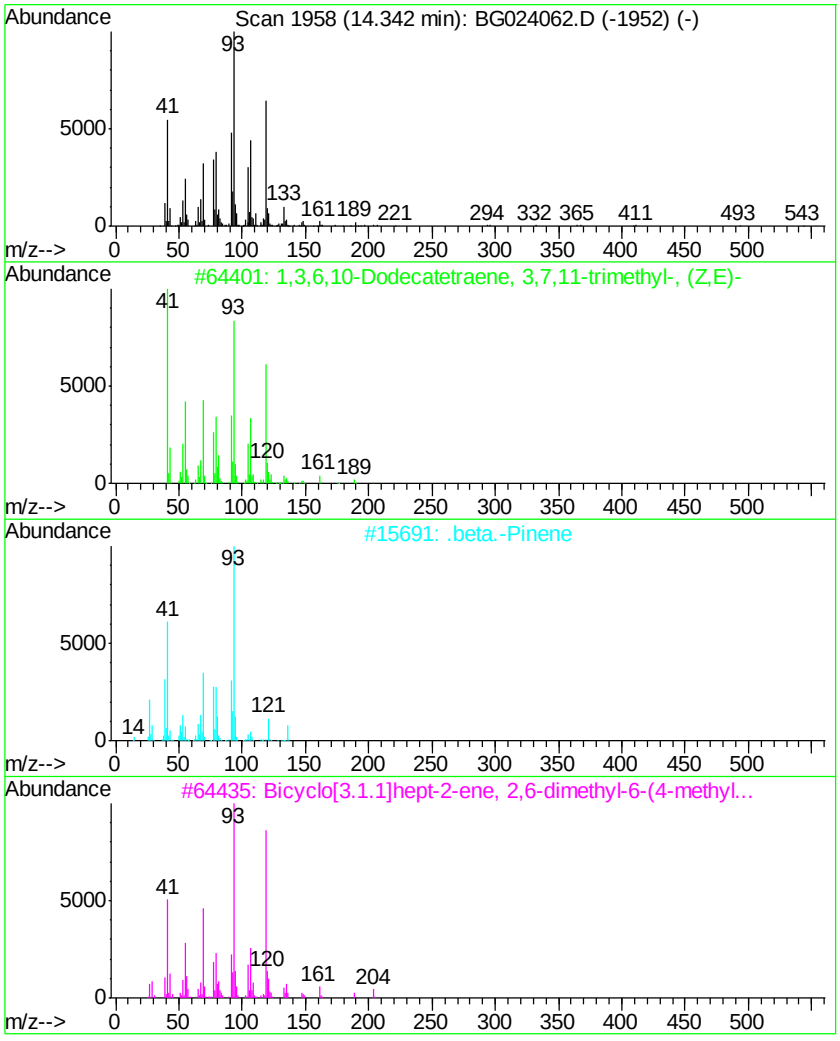
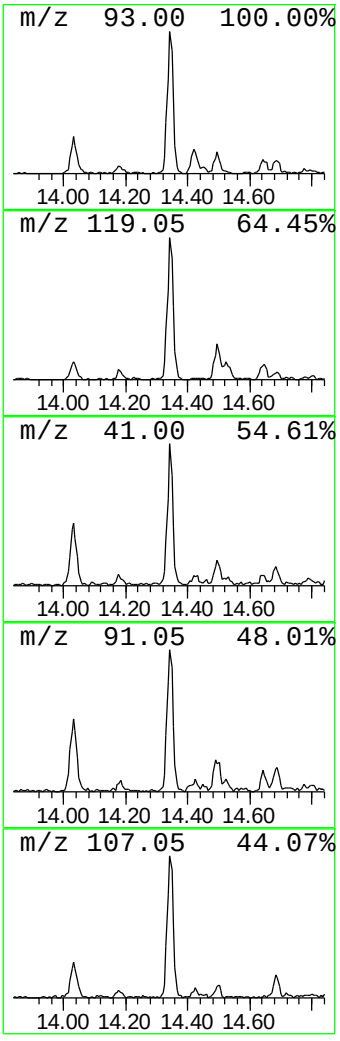
Peak Number 7 1,3,6,10-Dodecatetraene, 3,... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.34	10.88 ng	654520	Acenaphthene-d10	14.73

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,3,6,10-Dodecatetraene, 3,7,11-...	204	C15H24	026560-14-5	94
2	.beta.-Pinene	136	C10H16	000127-91-3	53
3	Bicyclo[3.1.1]hept-2-ene, 2,6-di...	204	C15H24	017699-05-7	50
4	1,3,7-Octatriene, 3,7-dimethyl-	136	C10H16	000502-99-8	50
5	1,3,6-Heptatriene, 2,5,5-trimethyl-	136	C10H16	029548-02-5	50





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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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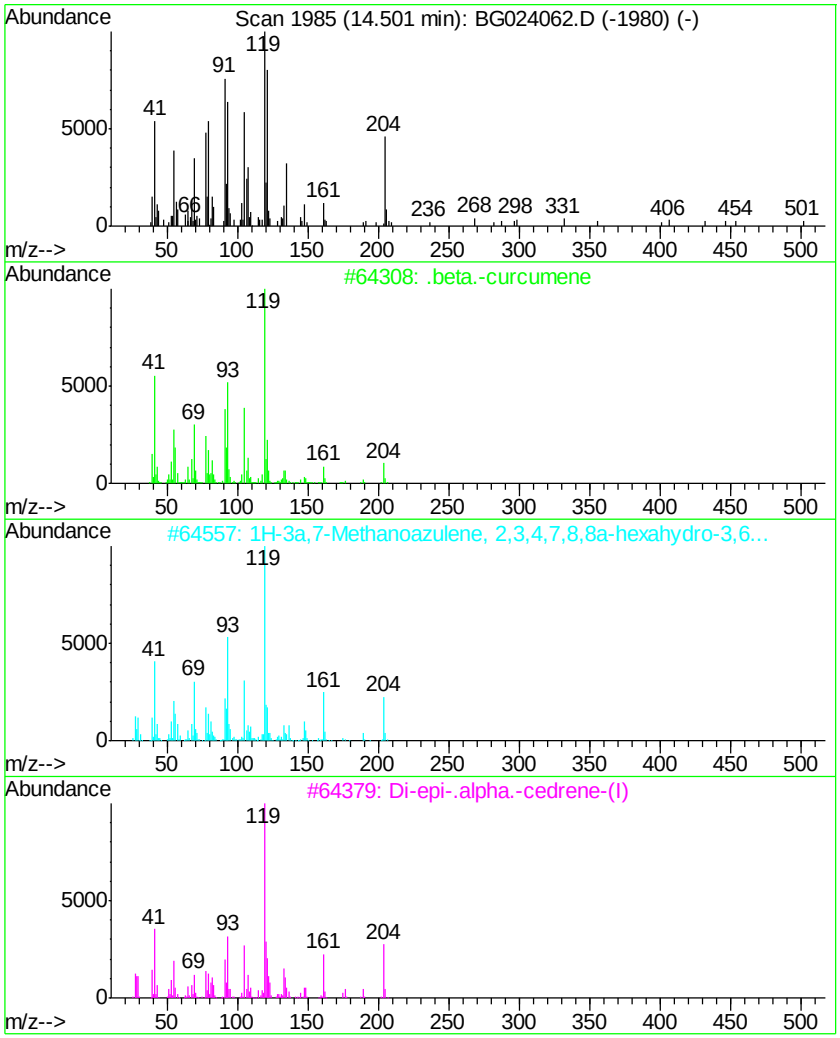
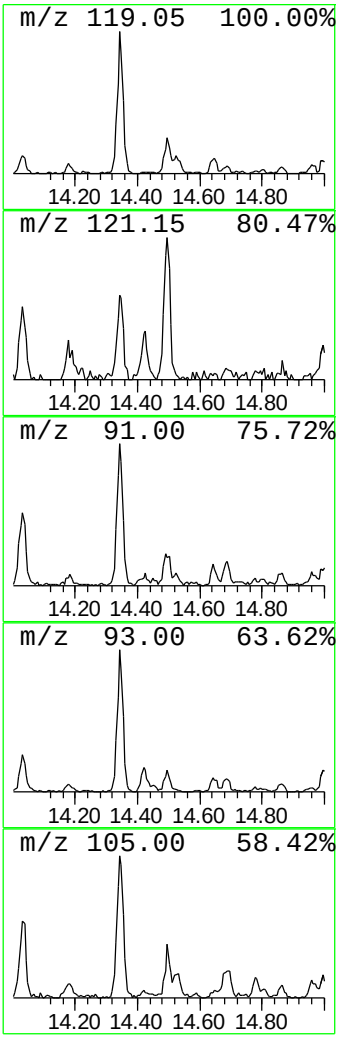
Peak Number 8 .beta.-curcumene Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.50	2.56 ng	154222	Acenaphthene-d10	14.73

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	.beta.-curcumene	204	C15H24	1000374-17-4	62
2	1H-3a,7-Methanoazulene, 2,3,4,7,...	204	C15H24	000469-61-4	60
3	Di-epi-.alpha.-cedrene-(I)	204	C15H24	021996-77-0	55
4	1H-Benzocycloheptene, 2,4a,5,6,7...	204	C15H24	001461-03-6	49
5	Bicyclo[3.1.1]hept-2-ene, 2,6-di...	204	C15H24	017699-05-7	49





Data Path : Z:\HPCHEM1\BNA G\DATA\BG092116\
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Sample : H4821-13 5X
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ALS Vial : 12 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : YORK-SB1-4

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

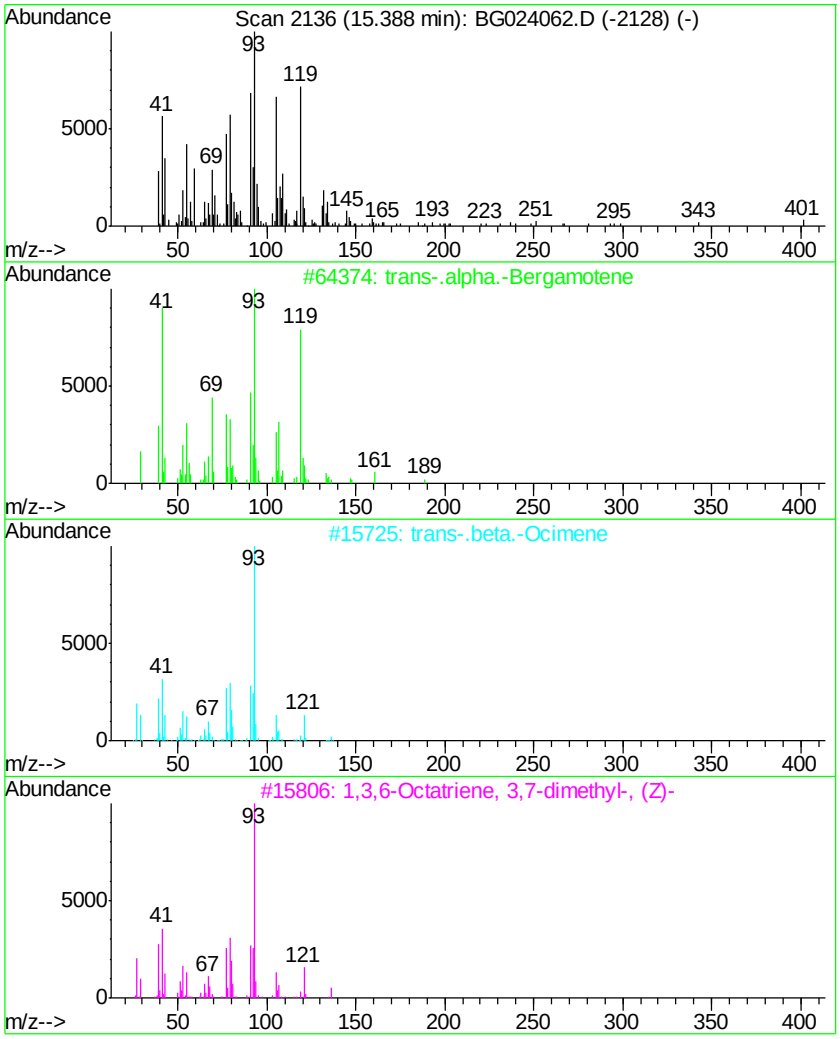
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

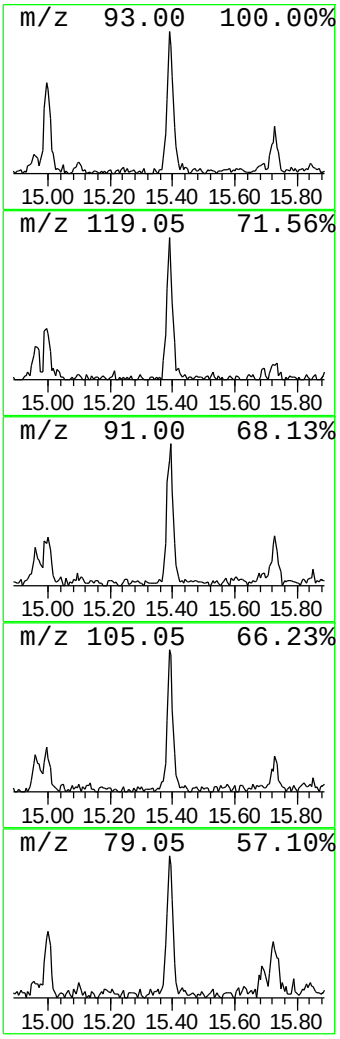
****

Peak Number 9 trans-.alpha.-Bergamotene Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.39	4.12 ng	247616	Acenaphthene-d10	14.73

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	trans-.alpha.-Bergamotene	204	C15H24	013474-59-4	50
2	trans-.beta.-Ocimene	136	C10H16	003779-61-1	38
3	1,3,6-Octatriene, 3,7-dimethyl-,...	136	C10H16	003338-55-4	38
4	.alpha.-Pinene	136	C10H16	000080-56-8	38
5	3-Methylenecycloheptene	108	C8H12	034564-56-2	30





Data Path : Z:\HPCHEM1\BNA G\DATA\BG092116\
Data File : BG024062.D
Acq On : 21 Sep 2016 19:37
Operator : UM/SJ
Sample : H4821-13 5X
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : YORK-SB1-4

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

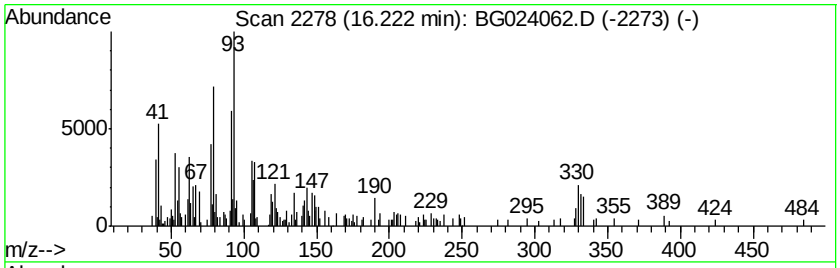
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

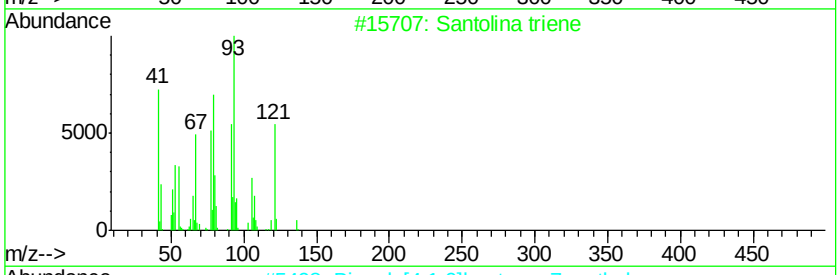
****

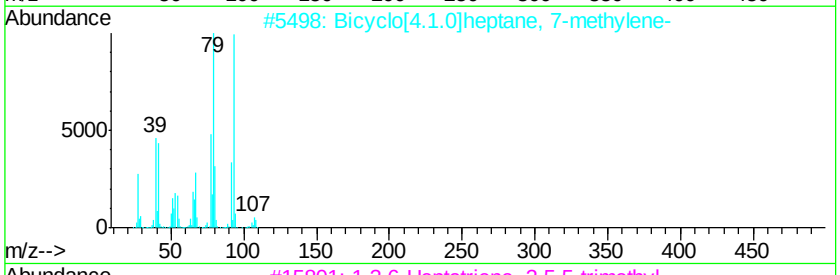
Peak Number 10 Santolina triene Concentration Rank 11

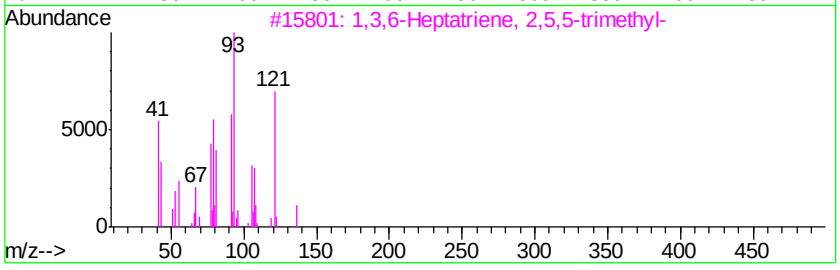
R.T.	EstConc	Area	Relative to ISTD	R.T.
16.22	2.76 ng	178480	Phenanthrene-d10	17.49

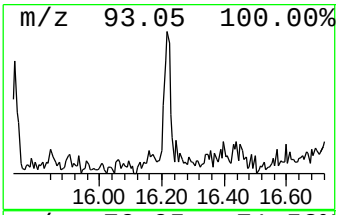
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Santolina triene	136	C10H16	002153-66-4	64
2	Bicyclo[4.1.0]heptane, 7-methylene-	108	C8H12	054211-14-2	60
3	1,3,6-Heptatriene, 2,5,5-trimethyl-	136	C10H16	029548-02-5	58
4	1,5,9,11-Tridecatetraene, 12-met...	190	C14H22	062338-27-6	58
5	Bicyclo[2.2.1]heptane, 2,2-dimet...	136	C10H16	005794-03-6	52

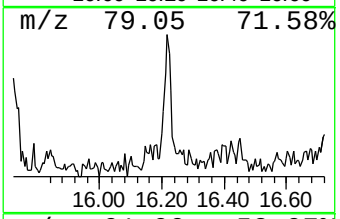


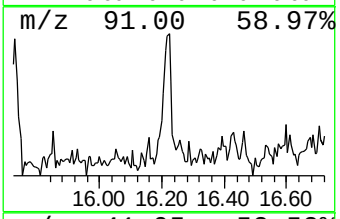


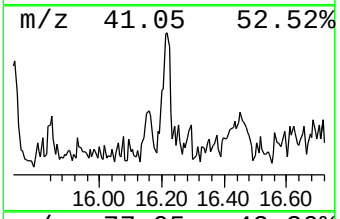


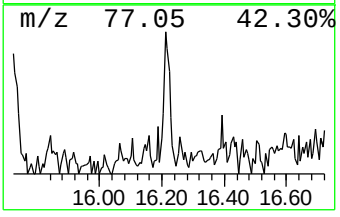












Data Path : Z:\HPCHEM1\BNA G\DATA\BG092116\
Data File : BG024062.D
Acq On : 21 Sep 2016 19:37
Operator : UM/SJ
Sample : H4821-13 5X
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : YORK-SB1-4

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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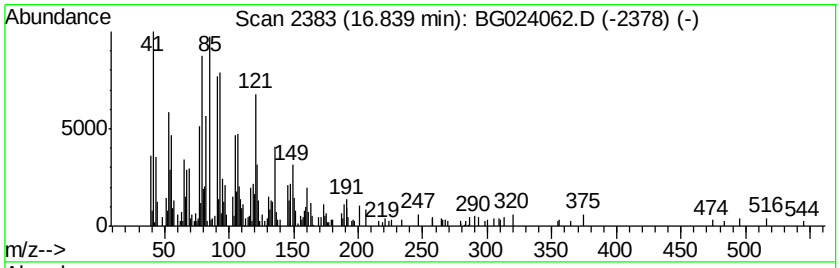
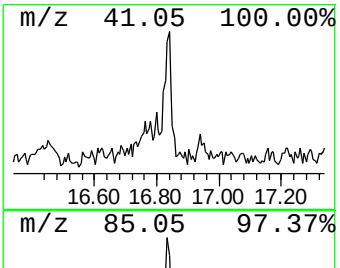
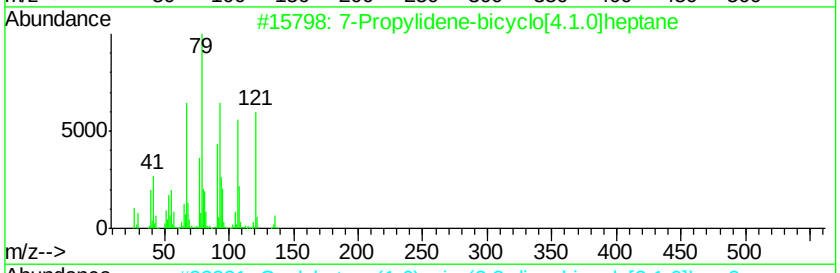
Peak Number 11 unknown16.84 Concentration Rank 9

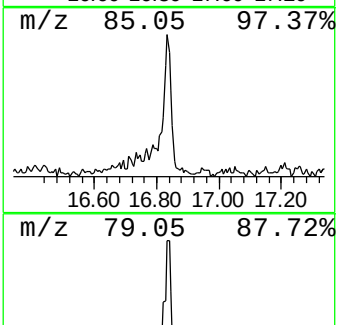
R.T.	EstConc	Area	Relative to ISTD	R.T.
16.84	3.93 ng	254169	Phenanthrene-d10	17.49

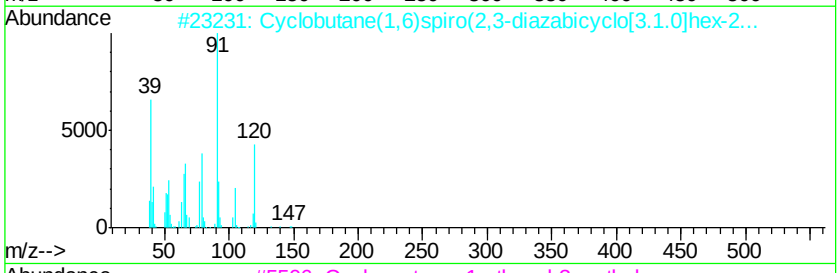
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	7-Propylidene-bicyclo[4.1.0]heptane	136	C10H16	082253-09-6	42
2	Cyclobutane(1,6)spiro(2,3-diazab...	148	C9H12N2	176172-15-9	38
3	Cyclopentane, 1-ethenyl-3-methyl...	108	C8H12	029668-63-1	35
4	7-Pentadecen-5-yne, (Z)-	206	C15H26	074744-49-3	25
5	1-Methylene-2-vinylcyclopentane	108	C8H12	006196-78-7	25

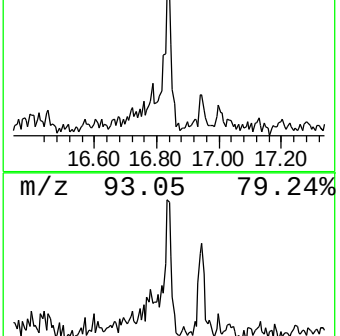


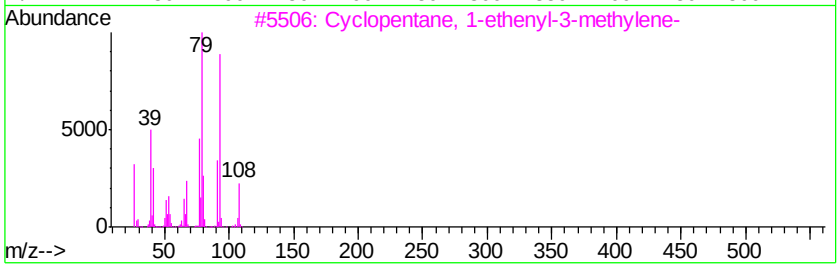


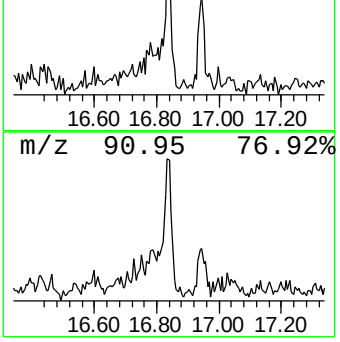












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Operator : UM/SJ
Sample : H4821-13 5X
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : YORK-SB1-4

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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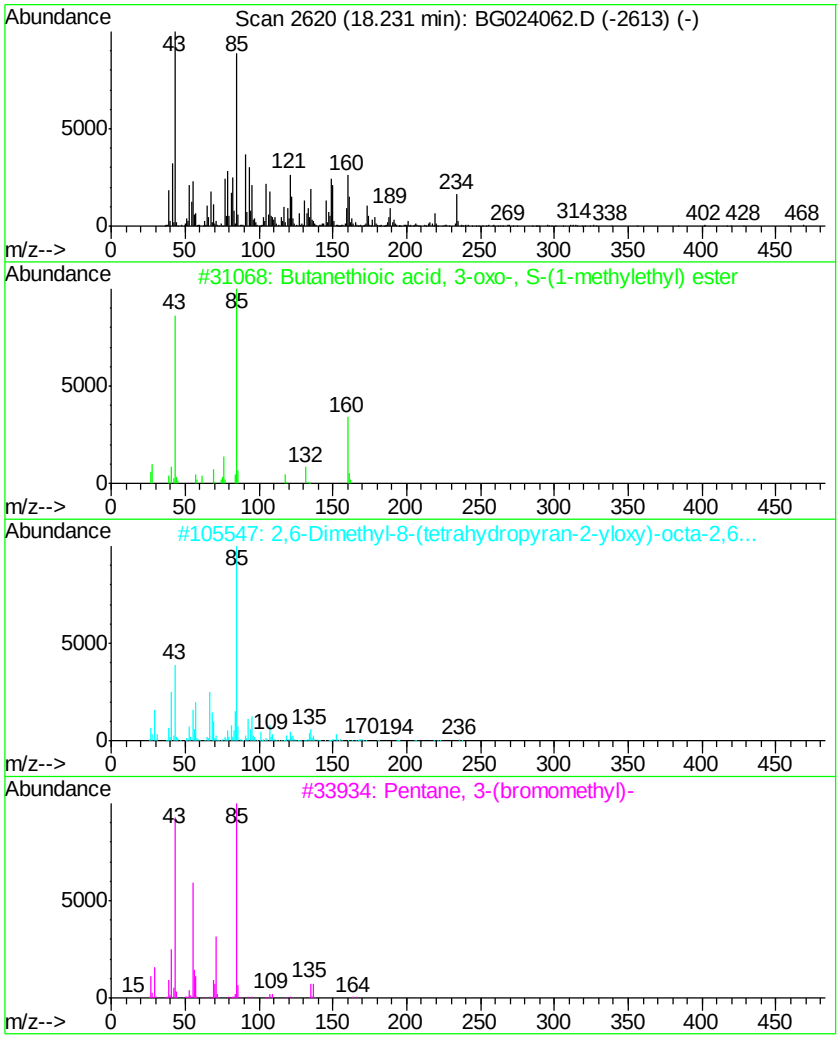
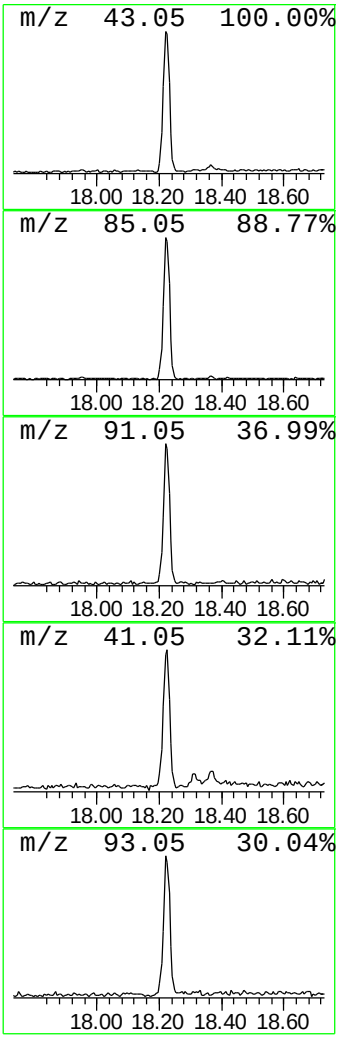
Peak Number 12 unknown18.23 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.23	25.39 ng	1640280	Phenanthrene-d10	17.49

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butanethioic acid, 3-oxo-, S-(1-...	160	C7H12O2S	015780-62-8	35
2	2,6-Dimethyl-8-(tetrahydropyran-...	254	C15H26O3	038290-53-8	27
3	Pentane, 3-(bromomethyl)-	164	C6H13Br	003814-34-4	25
4	2(3H)-Furanone, 5-butyldihydro-	142	C8H14O2	000104-50-7	22
5	2H-Pyran, 2-(8-dodecynyloxy)tetr...	266	C17H30O2	064604-68-8	22





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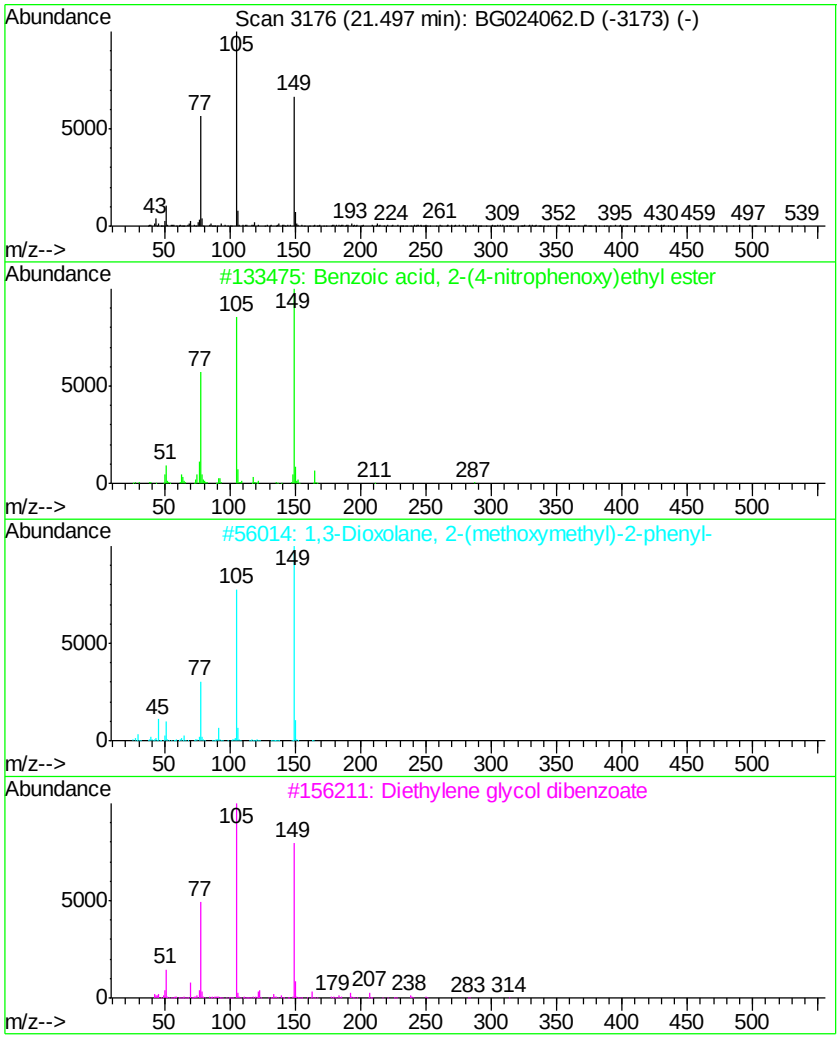
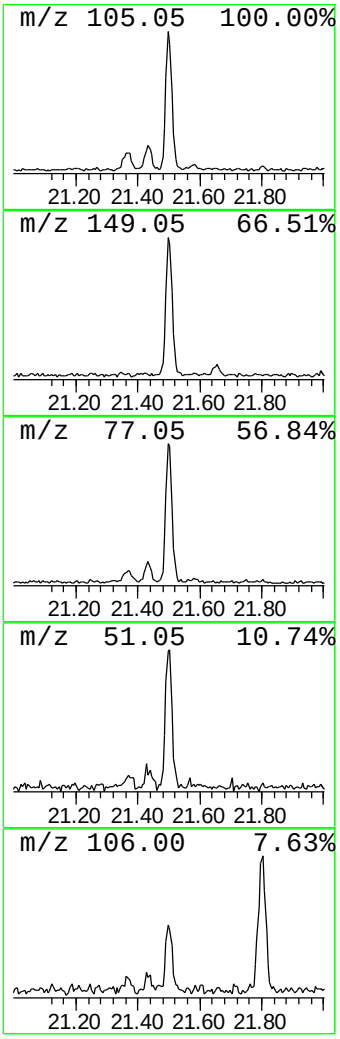
Peak Number 13 Benzoic acid, 2-(4-nitrophe... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.50	5.34 ng	399329	Chrysene-d12	21.80

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzoic acid, 2-(4-nitrophenoxy)...	287	C15H13NO5	1000368-92-7	59
2	1,3-Dioxolane, 2-(methoxymethyl)...	194	C11H14O3	1000156-71-2	50
3	Diethylene glycol dibenzoate	314	C18H18O5	000120-55-8	47
4	Quinazolin-4(3H)-one, 2-(1-methy...	324	C18H16N2O2S	298686-75-6	38
5	Benzenepropanenitrile, .beta.-oxo-	145	C9H7NO	000614-16-4	38





Tentatively Identified Compound (LSC) summary

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Data File : BG024062.D
Acq On : 21 Sep 2016 19:37
Operator : UM/SJ
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Misc :
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Instrument : BNA_G
ClientSampleId : YORK-SB1-4

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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...	2.93	25.0	ng	822405	1	8.05	658726	20.0
2-Pentanone, 4-hy...	5.13	3.2	ng	106374	1	8.05	658726	20.0
unknown7.59	7.59	6.6	ng	217559	1	8.05	658726	20.0
Bicyclo[7.2.0]und...	14.03	6.3	ng	376169	3	14.73	1203050	20.0
Dimethyl phthalate	14.18	2.1	ng	128497	3	14.73	1203050	20.0
1,3,6,10-Dodecate...	14.34	10.9	ng	654520	3	14.73	1203050	20.0
.beta.-curcumene	14.50	2.6	ng	154222	3	14.73	1203050	20.0
trans-.alpha.-Ber...	15.39	4.1	ng	247616	3	14.73	1203050	20.0
Santolina triene	16.22	2.8	ng	178480	4	17.49	1291910	20.0
unknown16.84	16.84	3.9	ng	254169	4	17.49	1291910	20.0
unknown18.23	18.23	25.4	ng	1640280	4	17.49	1291910	20.0
Benzoic acid, 2-(...	21.50	5.3	ng	399329	5	21.80	1495080	20.0