LSC Area Percent Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG031519\
Data File : BG039947.D
Acq On : 15 Mar 2019 18:49
Operator : JU/SJ
Sample : K1909-01
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : MH-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:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG030419.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	3.197	13	18	27	rBV	290325	578671	22.33%	3.349%
2	3.273	27	31	50	rVB	478101	723621	27.93%	4.188%
3	5.230	359	364	374	rBV2	19421	31445	1.21%	0.182%
4	5.688	436	442	452	rVB	194184	314002	12.12%	1.817%
5	7.292	709	715	726	rVB	122413	225462	8.70%	1.305%
6	7.673	771	780	794	rBV	390179	673236	25.98%	3.897%
7	8.149	855	861	868	rVB	159985	277582	10.71%	1.607%
8	8.472	907	916	924	rVB	579122	1010090	38.98%	5.846%
9	9.324	1052	1061	1069	rBV	501925	881810	34.03%	5.104%
10	9.864	1143	1153	1164	rVB3	39934	68146	2.63%	0.394%
11	10.376	1231	1240	1244	rBV3	17559	31114	1.20%	0.180%
12	10.728	1292	1300	1308	rBV3	34191	66363	2.56%	0.384%
13	10.816	1308	1315	1319	rBV3	27236	51050	1.97%	0.295%
14	10.969	1334	1341	1345	rBV	225387	436660	16.85%	2.527%
15	11.004	1345	1347	1354	rVB	144718	202578	7.82%	1.173%
16	12.672	1624	1631	1637	rBV3	21785	36108	1.39%	0.209%
17	13.395	1745	1754	1761	rBV	1335799	2099787	81.04%	12.154%
18	14.217	1886	1894	1898	rBV	34872	53157	2.05%	0.308%
19	14.682	1968	1973	1981	rBV	15452	27038	1.04%	0.156%
20	14.775	1981	1989	1998	rVV2	356151	592474	22.87%	3.429%
21	15.281	2066	2075	2087	rBV	293923	548357	21.16%	3.174%
22	15.410	2091	2097	2103	rVB6	20417	36494	1.41%	0.211%
23	15.809	2161	2165	2172	rBV7	17168	31566	1.22%	0.183%
24	16.256	2235	2241	2251	rBV	626660	997295	38.49%	5.772%
25	17.519	2449	2456	2469	rBV2	479666	789770	30.48%	4.571%
26	18.095	2548	2554	2558	rBV7	20339	36028	1.39%	0.209%
27	18.265	2576	2583	2594	rBV5	56962	151345	5.84%	0.876%
28	18.371	2597	2601	2605	rBV2	41195	56778	2.19%	0.329%
29	18.517	2620	2626	2639	rBV	479965	768908	29.68%	4.450%
30	18.658	2647	2650	2657	rVB3	38830	61910	2.39%	0.358%
31	18.729	2657	2662	2671	rVV	70105	127638	4.93%	0.739%
32	18.805	2671	2675	2681	rVB2	48537	65749	2.54%	0.381%
33	18.987	2703	2706	2711	rVB6	21294	28937	1.12%	0.167%
34	19.105	2721	2726	2734	rBV	102300	172135	6.64%	0.996%

LSC Area Percent Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG031519\
Data File : BG039947.D
Acq On : 15 Mar 2019 18:49
Operator : JU/SJ
Sample : K1909-01
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : MH-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:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG030419.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

35	19.322	2760	2763	2771	rVB3	33173	51590	1.99%	0.299%
36	19.634	2812	2816	2823	rVB6	20377	33447	1.29%	0.194%
37	20.109	2891	2897	2903	rBV	1881388	2590973	100.00%	14.997%
38	20.650	2985	2989	2997	rVB10	12334	27769	1.07%	0.161%
39	20.879	3024	3028	3033	rBV7	26764	43244	1.67%	0.250%
40	21.402	3112	3117	3123	rBV10	24997	55161	2.13%	0.319%
41	21.807	3180	3186	3193	rVB2	511409	855170	33.01%	4.950%
42	22.036	3221	3225	3230	rVB7	17278	27783	1.07%	0.161%
43	22.571	3312	3316	3323	rVB3	21216	32998	1.27%	0.191%
44	22.759	3342	3348	3356	rBV2	49328	102969	3.97%	0.596%
45	23.288	3433	3438	3444	rVB3	31823	54253	2.09%	0.314%
46	24.122	3573	3580	3588	rVB3	40919	90510	3.49%	0.524%
47	25.114	3740	3749	3750	rBV4	34946	76797	2.96%	0.445%
48	25.173	3750	3759	3771	rVB2	289698	867430	33.48%	5.021%
49	26.283	3940	3948	3957	rVB3	26533	83011	3.20%	0.480%
50	27.693	4180	4188	4190	rBV8	13419	30749	1.19%	0.178%

Sum of corrected areas: 17277158

LSC Report - Integrated Chromatogram

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Data File : BG039947.D
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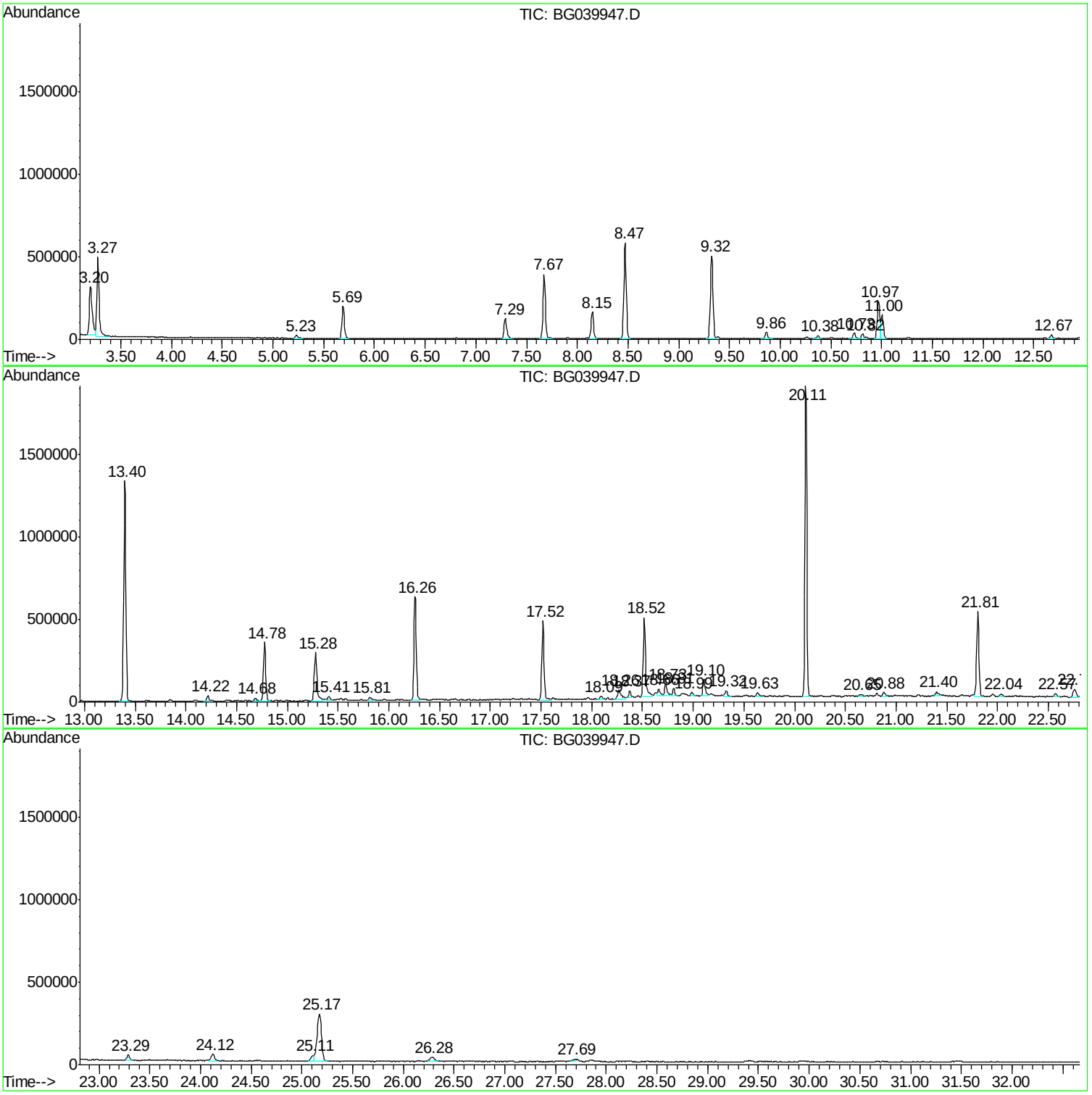
Operator : JU/SJ
Sample : K1909-01
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : MH-02

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG030419.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:\SVOASRV\HPCHEM1\BNA G\DATA\BG031519\
Data File : BG039947.D
Acq On : 15 Mar 2019 18:49
Operator : JU/SJ
Sample : K1909-01
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : MH-02

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG030419.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

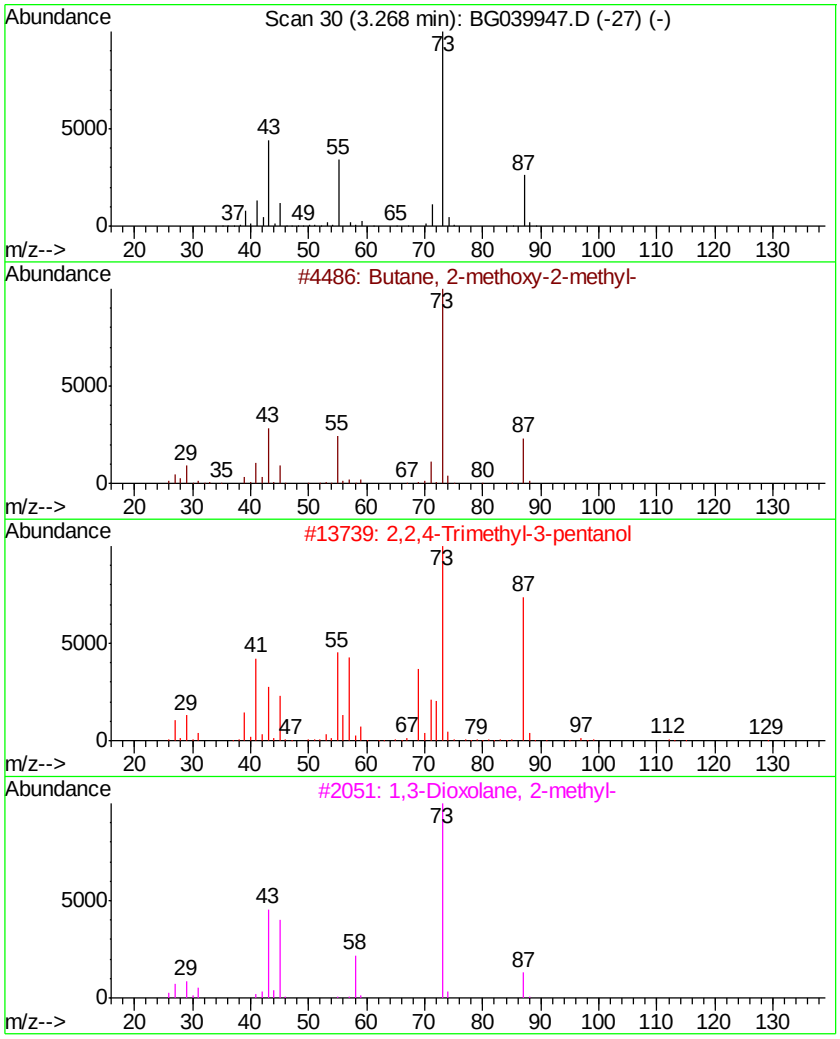
TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

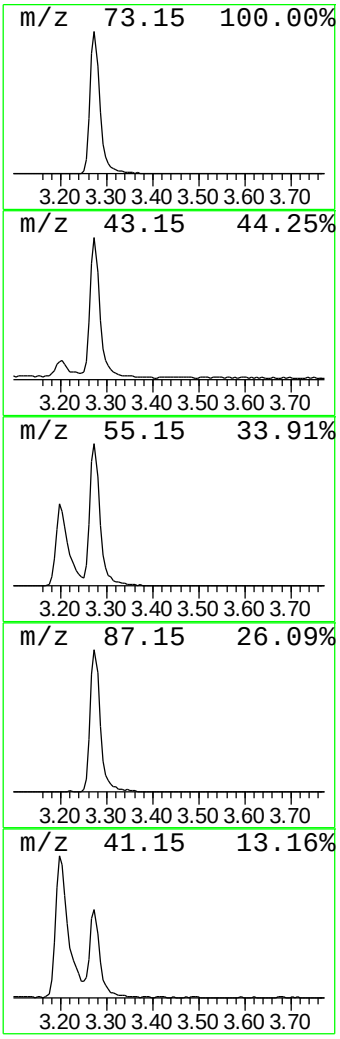
****

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.27	52.14 ng	723621	1,4-Dichlorobenzene-d4	8.15

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	83
2	2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	39
3	1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	25
4	Silane, tetramethyl-	88	C4H12Si	000075-76-3	12
5	Pentane, 3-methoxy-	102	C6H14O	036839-67-5	12





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Data File : BG039947.D
Acq On : 15 Mar 2019 18:49
Operator : JU/SJ
Sample : K1909-01
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : MH-02

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG030419.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

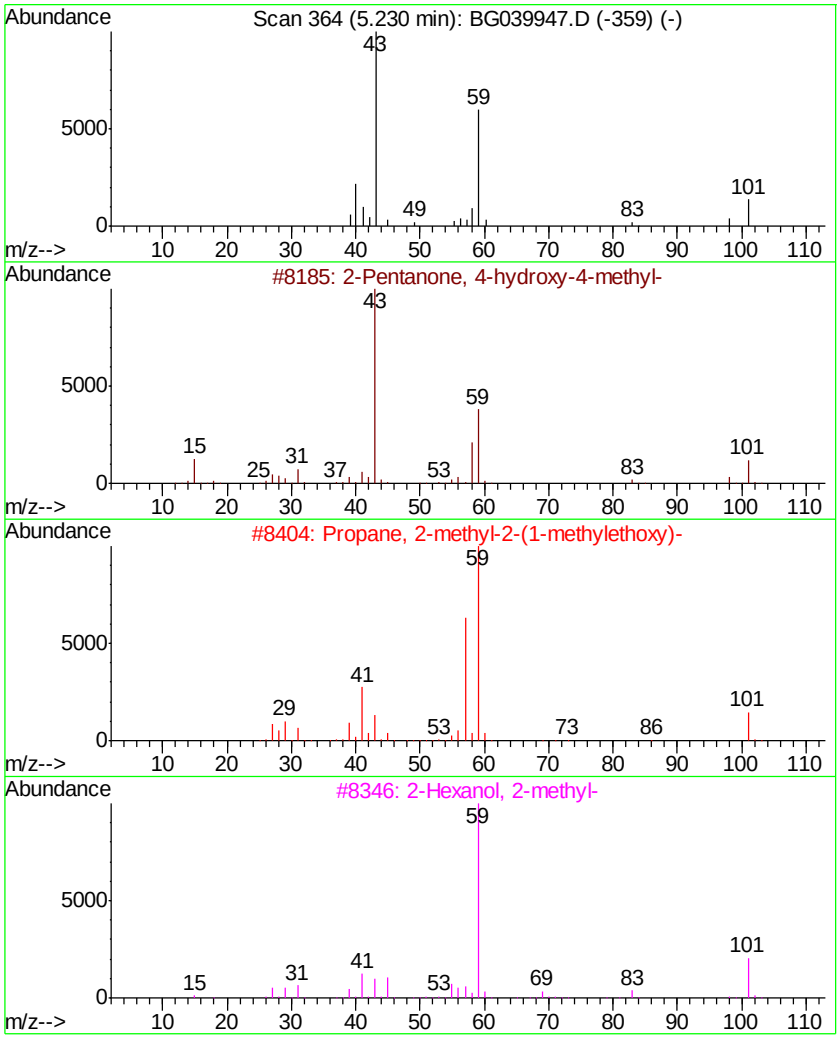
TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

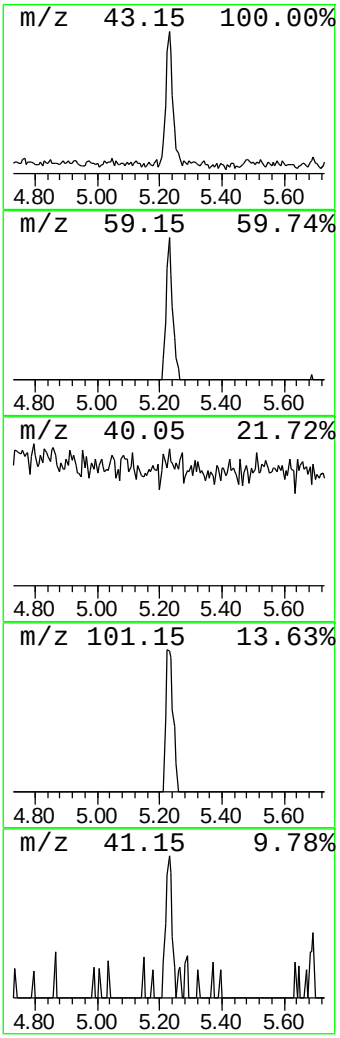
****

Peak Number 3 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.23	2.27 ng	31445	1,4-Dichlorobenzene-d4	8.15

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50
2	Propane, 2-methyl-2-(1-methyleth...	116	C7H16O	017348-59-3	42
3	2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	38
4	3-Octanol	130	C8H18O	000589-98-0	36
5	2-Pentanol, 2,4-dimethyl-	116	C7H16O	000625-06-9	33





Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG031519\
Data File : BG039947.D
Acq On : 15 Mar 2019 18:49
Operator : JU/SJ
Sample : K1909-01
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : MH-02

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG030419.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

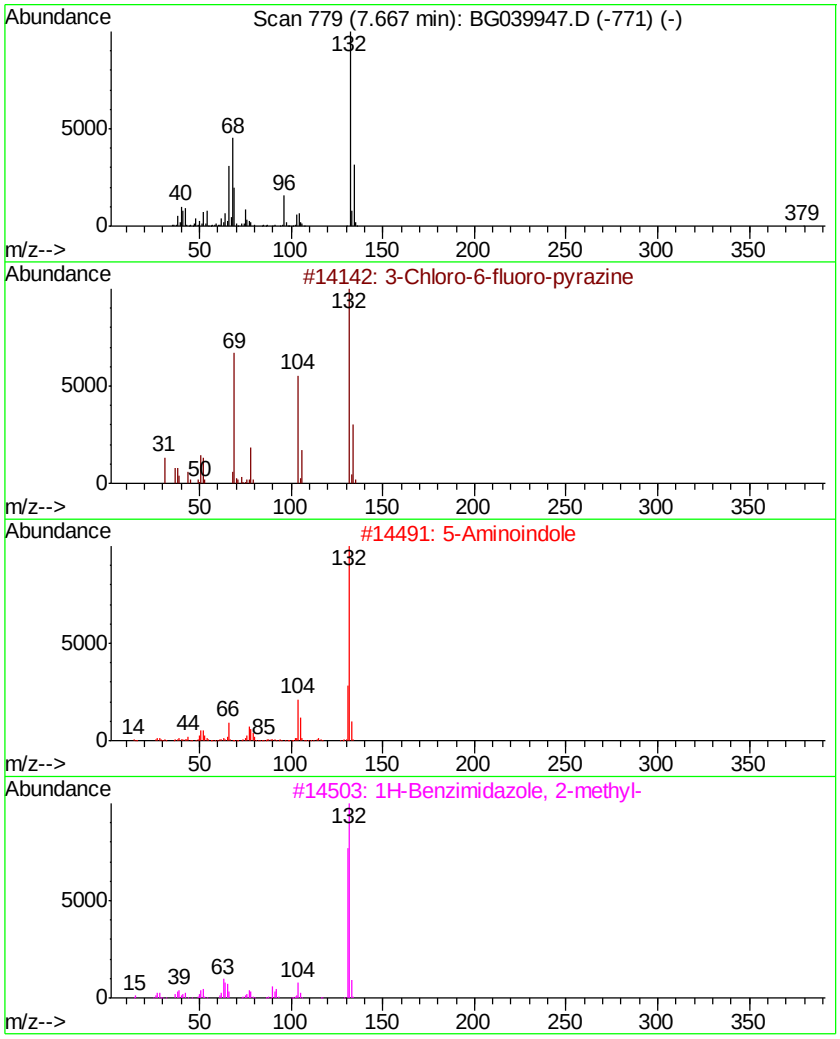
TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

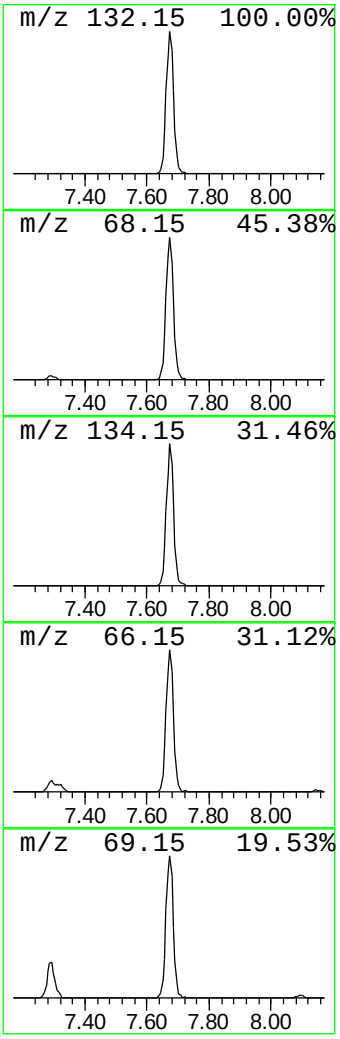
****

Peak Number 4 unknown7.67 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.67	48.51 ng	673236	1,4-Dichlorobenzene-d4	8.15

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	35
2	5-Aminoindole	132	C8H8N2	005192-03-0	12
3	1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	12
4	5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	12
5	N-(-3,4-Methylenedioxyphenylisop...	267	C17H17NO2	1000378-96-6	10





Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG031519\
Data File : BG039947.D
Acq On : 15 Mar 2019 18:49
Operator : JU/SJ
Sample : K1909-01
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : MH-02

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG030419.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

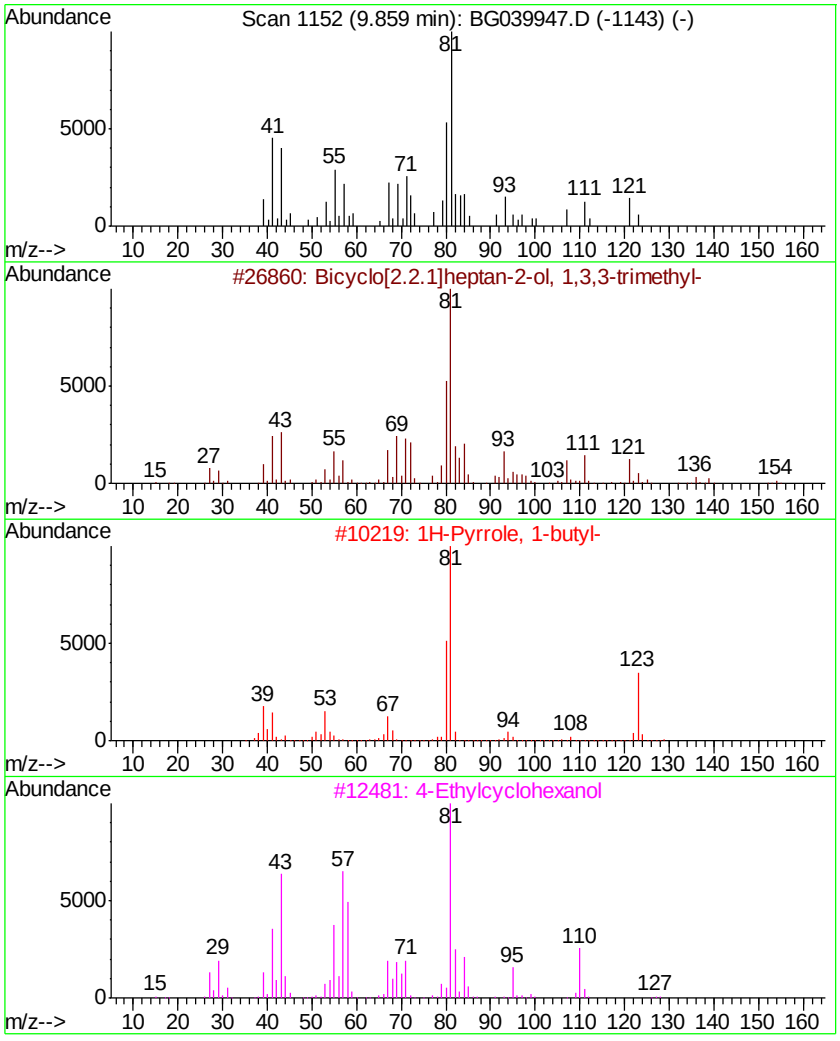
TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

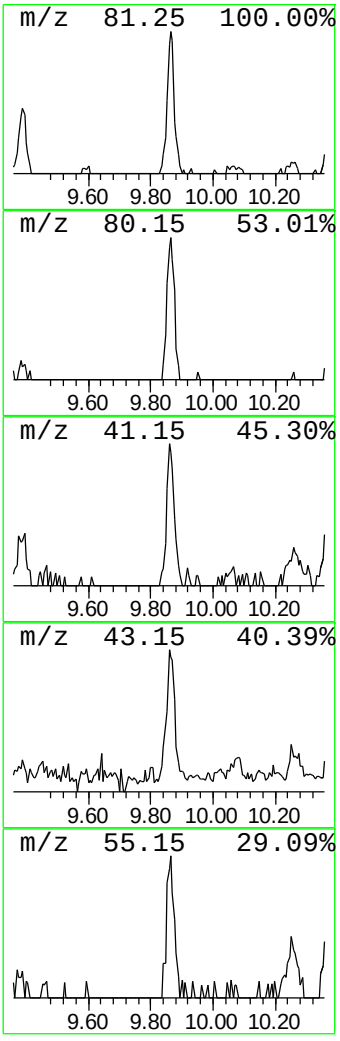
****

Peak Number 6 Bicyclo[2.2.1]heptan-2-ol, ... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.86	3.12 ng	68146	Naphthalene-d8	10.97

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Bicyclo[2.2.1]heptan-2-ol, 1,3,3...	154	C10H18O	001632-73-1	72
2	1H-Pyrrole, 1-butyl-	123	C8H13N	000589-33-3	43
3	4-Ethylcyclohexanol	128	C8H16O	004534-74-1	40
4	Cyclohexanol, 4-(1-methylethyl)-	142	C9H18O	004621-04-9	38
5	Fenchol, exo-	154	C10H18O	022627-95-8	38





Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG031519\
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Operator : JU/SJ
Sample : K1909-01
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : MH-02

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG030419.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

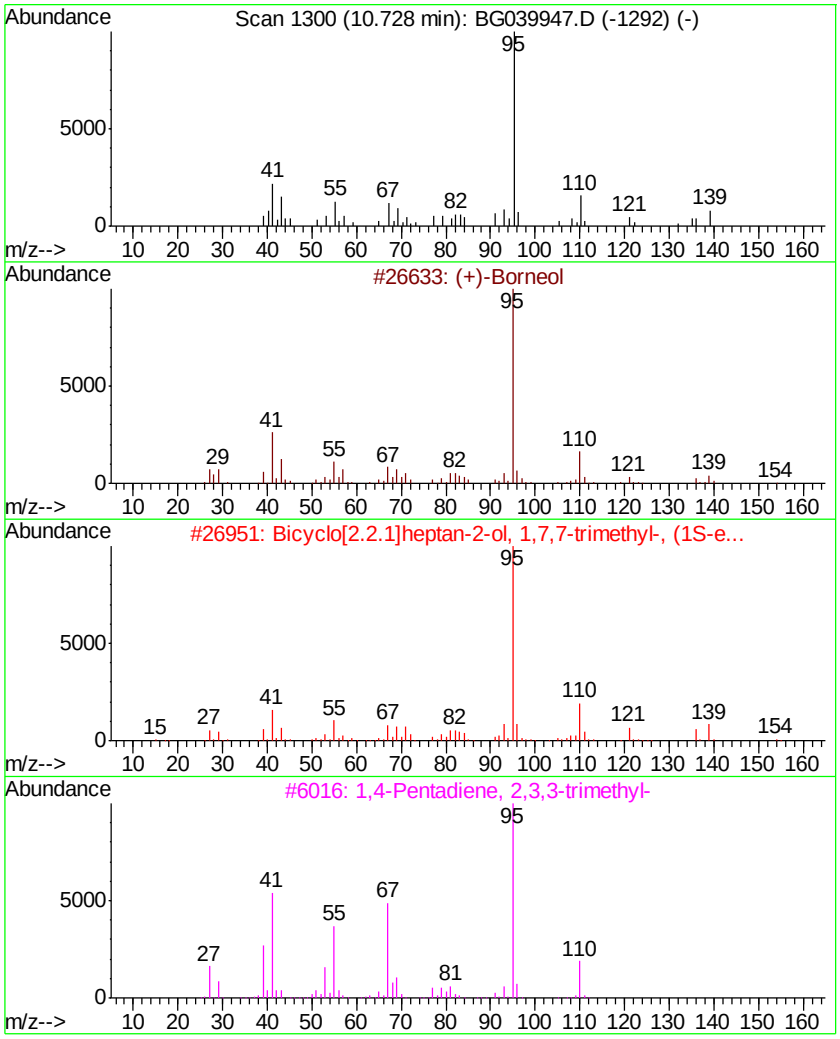
TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

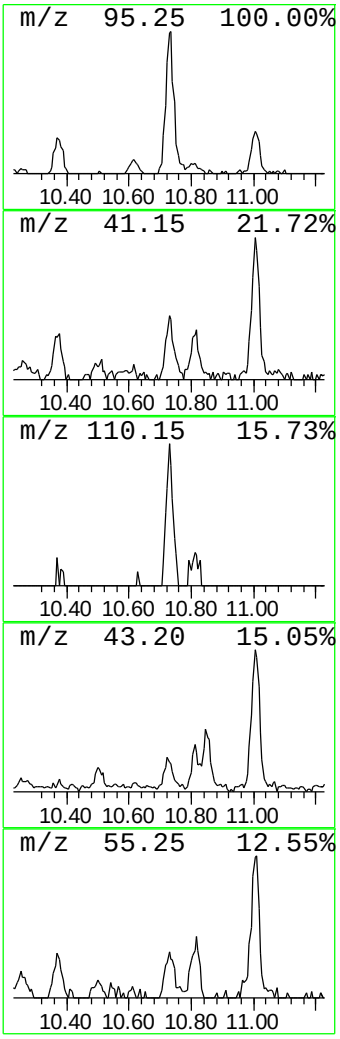
****

Peak Number 7 (+)-Borneol Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.73	3.04 ng	66363	Naphthalene-d8	10.97

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	(+)-Borneol	154	C10H18O	1000150-76-3	72
2	Bicyclo[2.2.1]heptan-2-ol, 1,7,7...	154	C10H18O	000464-45-9	64
3	1,4-Pentadiene, 2,3,3-trimethyl-	110	C8H14	000756-02-5	58
4	Furan, 2-ethyl-5-methyl-	110	C7H10O	001703-52-2	58
5	1,3-Dimethyl-1-cyclohexene	110	C8H14	002808-76-6	58





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Data File : BG039947.D
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Misc :
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Instrument : BNA_G
ClientSampleId : MH-02

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG030419.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

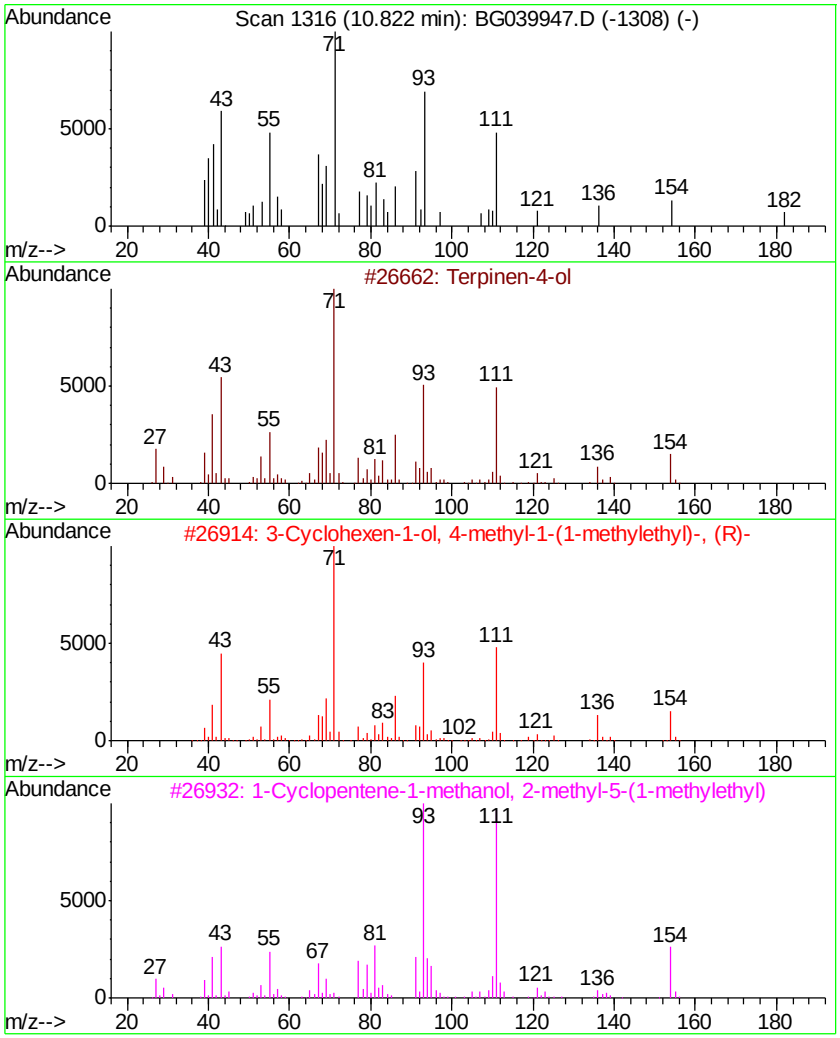
TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

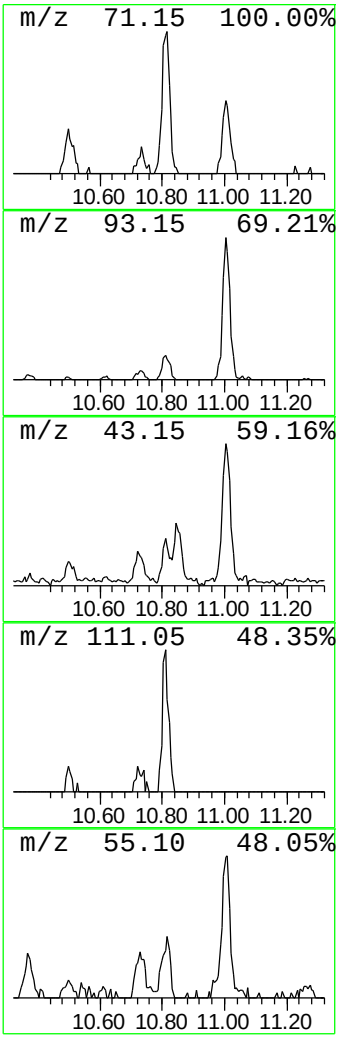
****

Peak Number 8 Terpinen-4-ol Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.82	2.34 ng	51050	Naphthalene-d8	10.97

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Terpinen-4-ol	154	C10H18O	000562-74-3	95
2	3-Cyclohexen-1-ol, 4-methyl-1-(1...	154	C10H18O	020126-76-5	68
3	1-Cyclopentene-1-methanol, 2-met...	154	C10H18O	080113-82-2	38
4	2-Buten-1-ol, 3-methyl-	86	C5H10O	000556-82-1	30
5	2-Octen-4-ol, (E)-	128	C8H16O	020125-81-9	27





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Misc :
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Instrument : BNA_G
ClientSampleId : MH-02

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG030419.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

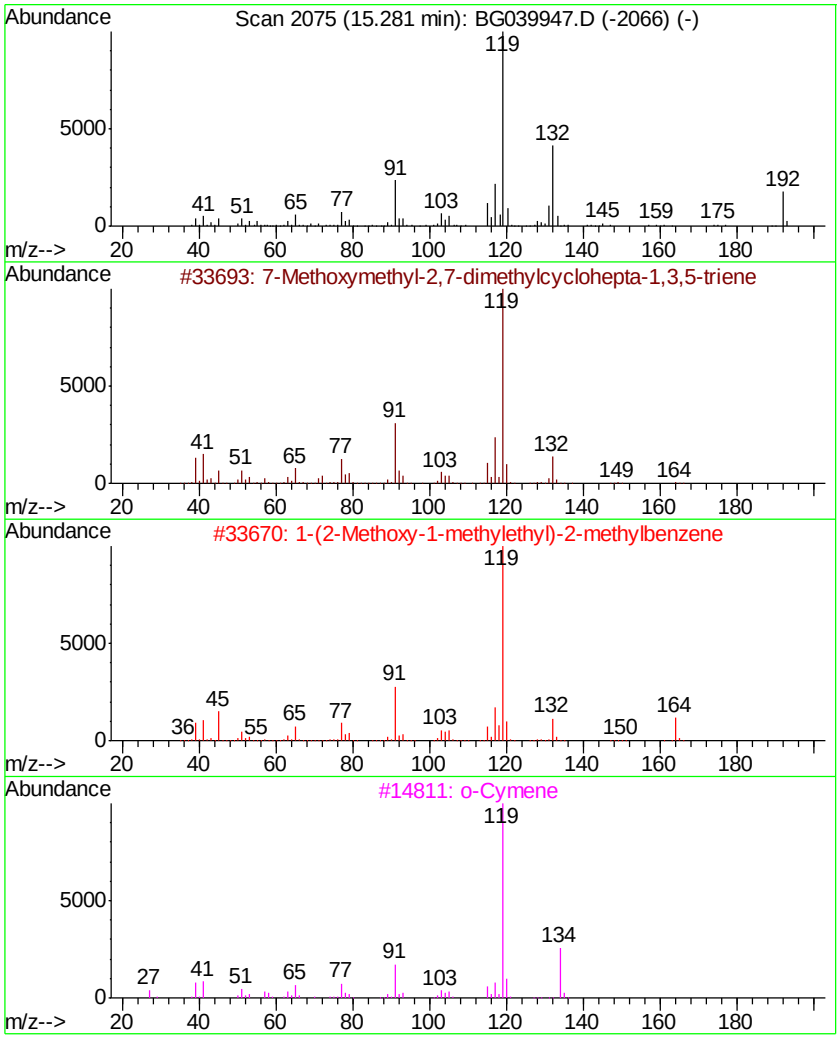
TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

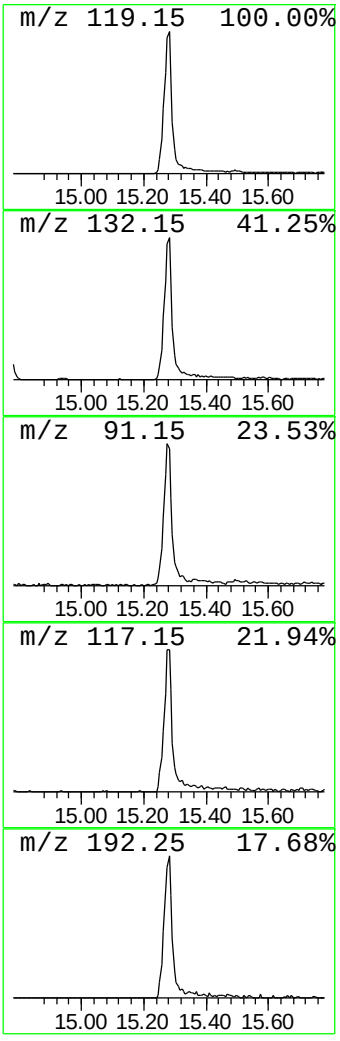
****

Peak Number 9 7-Methoxymethyl-2,7-dimethy... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.28	18.51 ng	548357	Acenaphthene-d10	14.78

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	7-Methoxymethyl-2,7-dimethylcycl...	164	C11H16O	073992-48-0	78
2	1-(2-Methoxy-1-methylethyl)-2-me...	164	C11H16O	1000210-02-1	47
3	o-Cymene	134	C10H14	000527-84-4	43
4	4-Oxo-4-(para-tolyl)-butyric acid	192	C11H12O3	004619-20-9	38
5	Ethyl 1,3-dithiane-2-carboxylate	192	C7H12O2S2	020462-00-4	38





Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG031519\
Data File : BG039947.D
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Sample : K1909-01
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : MH-02

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG030419.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

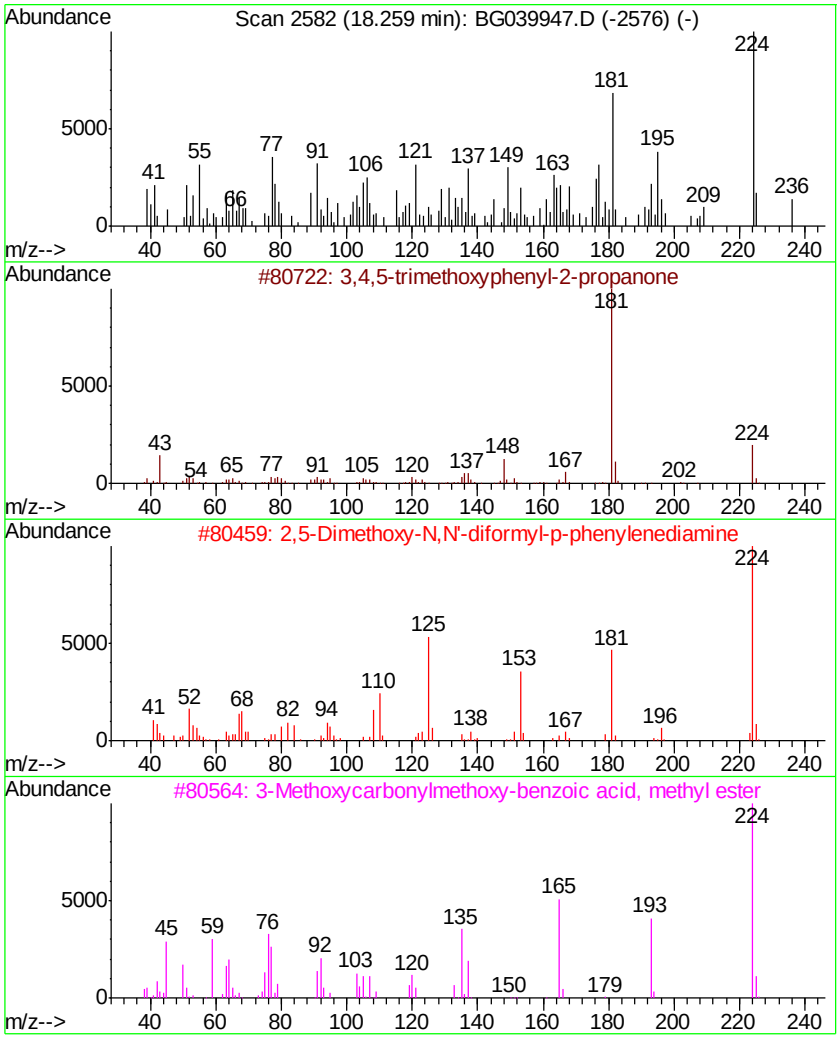
TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

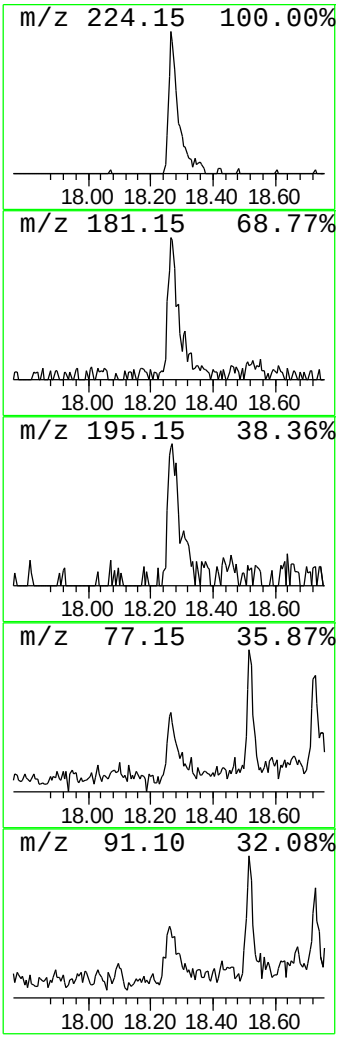
****

Peak Number 10 3,4,5-trimethoxyphenyl-2-pr... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.26	3.83 ng	151345	Phenanthrene-d10	17.52

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3,4,5-trimethoxyphenyl-2-propanone	224	C12H16O4	1000378-97-6	53
2	2,5-Dimethoxy-N,N'-diformyl-p-ph...	224	C10H12N2O4	1000250-95-2	52
3	3-Methoxycarbonylmethoxy-benzoic...	224	C11H12O5	1000317-51-5	50
4	Benzenebutyric acid, 2,3-dimethoxy-	224	C12H16O4	064400-76-6	49
5	1,1-Dimethyl-2-(diphenylmethylen...	224	C15H16N2	024398-55-8	46





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Data File : BG039947.D
Acq On : 15 Mar 2019 18:49
Operator : JU/SJ
Sample : K1909-01
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : MH-02

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG030419.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

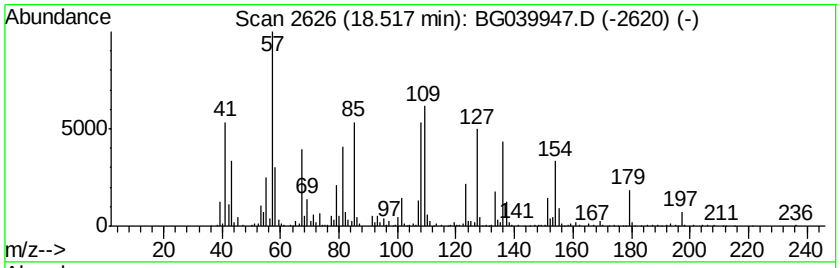
TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

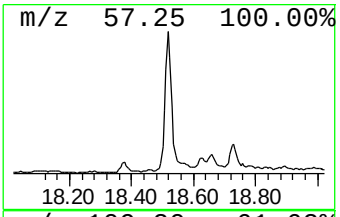
****

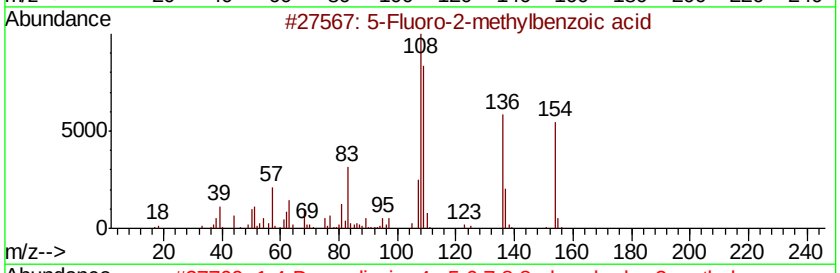
Peak Number 11 unknown18.52 Concentration Rank 5

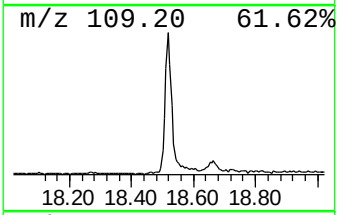
R.T.	EstConc	Area	Relative to ISTD	R.T.
18.52	19.47 ng	768908	Phenanthrene-d10	17.52

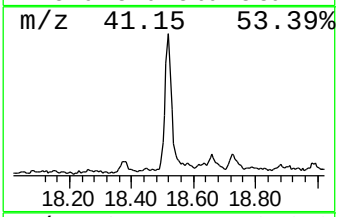
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	5-Fluoro-2-methylbenzoic acid	154	C8H7FO2	033184-16-6	46
2	1,4-Benzodioxin, 4a,5,6,7,8,8a-h...	154	C9H14O2	007196-96-5	25
3	Formic acid, (4-fluorophenyl)met...	154	C8H7FO2	1000368-72-9	22
4	1H-Pyrrole, 4-ethyl-2,3-dimethyl-	123	C8H13N	000491-18-9	22
5	2,5-Furandione, dihydro-3-(2-met...	154	C8H10O3	018908-20-8	22

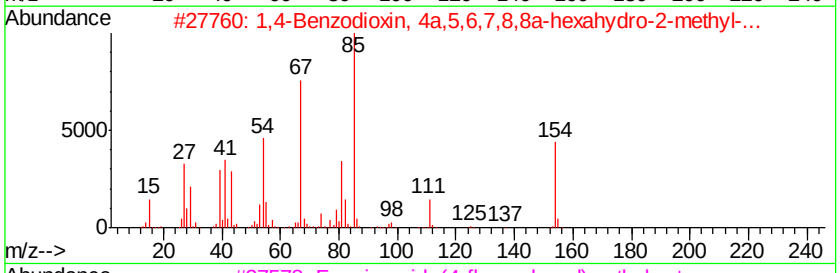


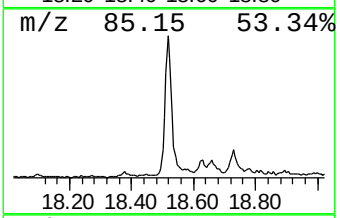


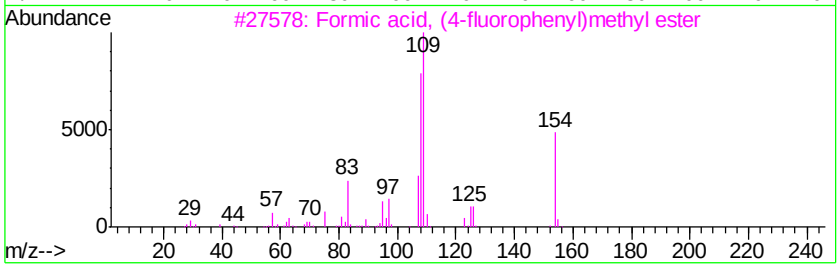


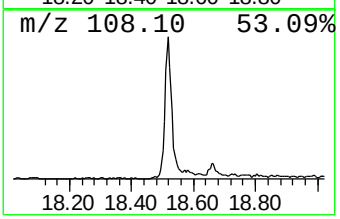












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Operator : JU/SJ
Sample : K1909-01
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : MH-02

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG030419.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

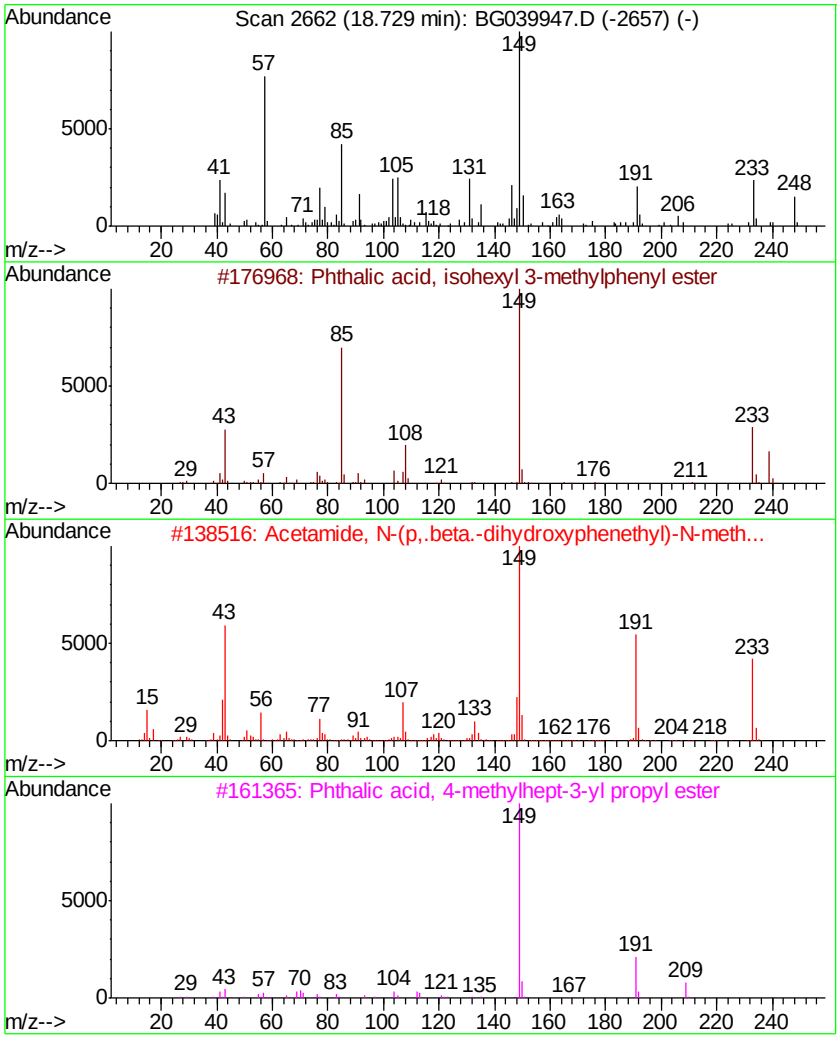
TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

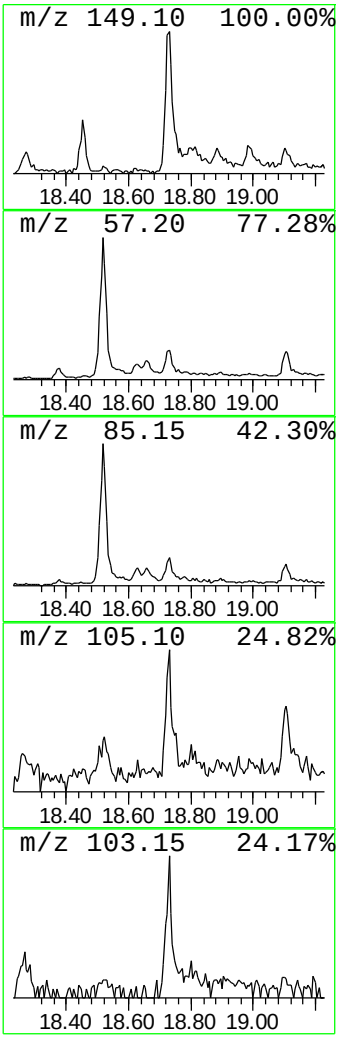
****

Peak Number 12 unknown18.73 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.73	3.23 ng	127638	Phenanthrene-d10	17.52

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phthalic acid, isohexyl 3-methyl...	340	C21H24O4	1000315-37-1	47
2	Acetamide, N-(p,.beta.-dihydroxy...	293	C15H19NO5	014383-57-4	35
3	Phthalic acid, 4-methylhept-3-yl...	320	C19H28O4	1000377-93-7	27
4	Phthalic acid, hexyl phenyl ester	326	C20H22O4	1000314-87-2	25
5	Phthalic acid, 2-cyclohexylethyl...	332	C20H28O4	1000309-06-9	22





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Instrument : BNA_G
ClientSampleId : MH-02

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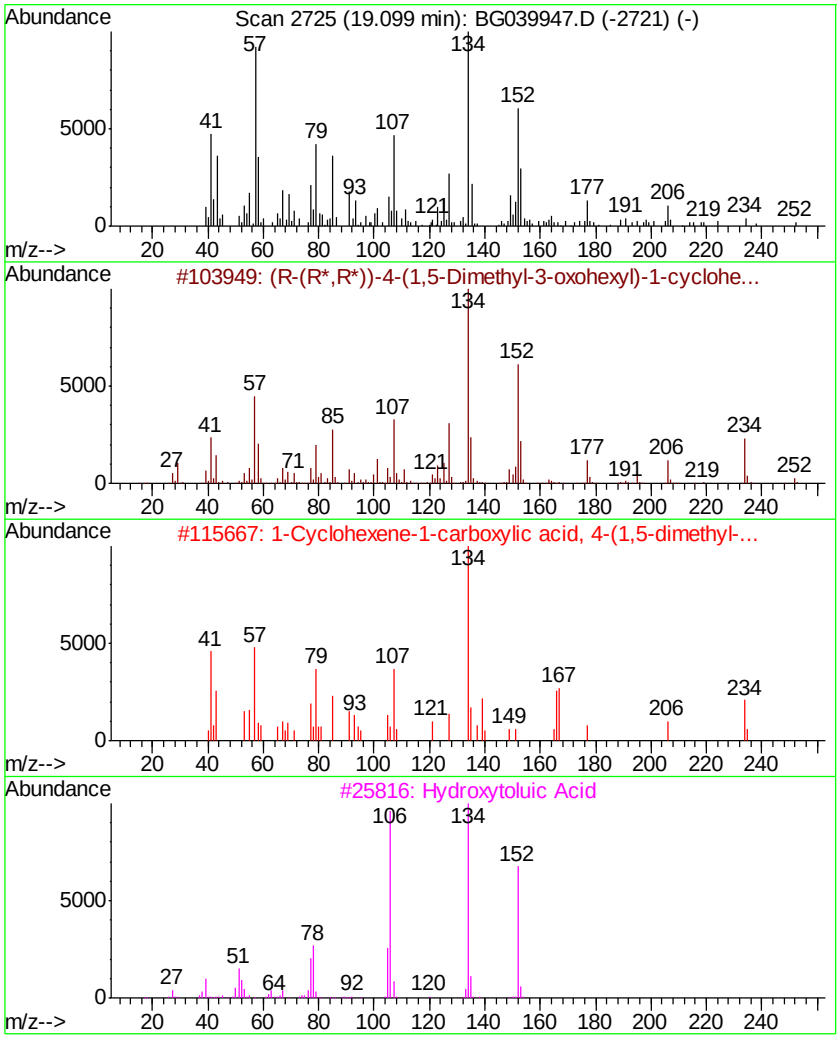
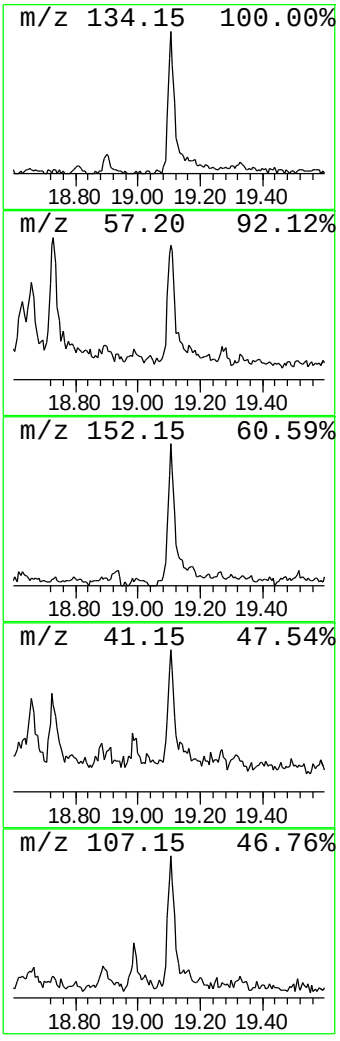
Peak Number 13 (R-(R*,R*))-4-(1,5-Dimethyl... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.10	4.36 ng	172135	Phenanthrene-d10	17.52

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	(R-(R*,R*))-4-(1,5-Dimethyl-3-ox...	252	C15H24O3	006753-22-6	74
2	1-Cyclohexene-1-carboxylic acid,...	266	C16H26O3	017904-27-7	53
3	Hydroxytoluic Acid	152	C8H8O3	000083-40-9	35
4	1H-Purine, 6-methyl-	134	C6H6N4	002004-03-7	27
5	2-Benzoxazolamine	134	C7H6N2O	004570-41-6	27





Library Search Compound Report

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Data File : BG039947.D
Acq On : 15 Mar 2019 18:49
Operator : JU/SJ
Sample : K1909-01
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : MH-02

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG030419.M
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TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

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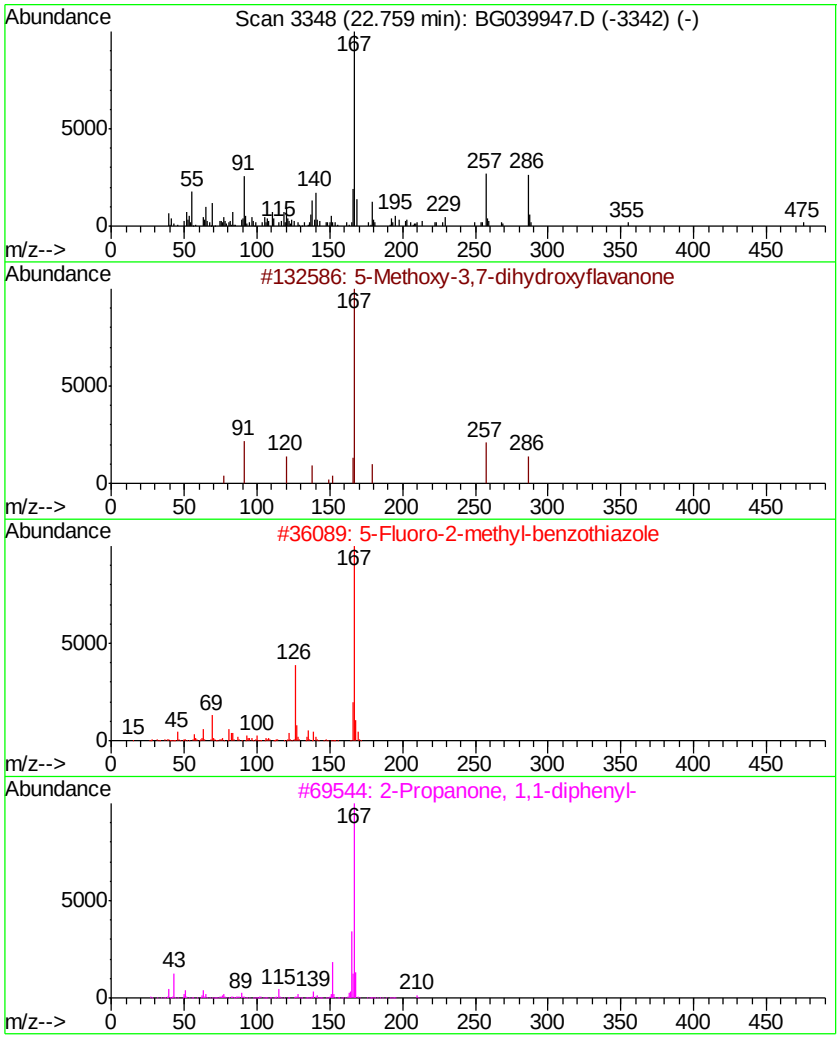
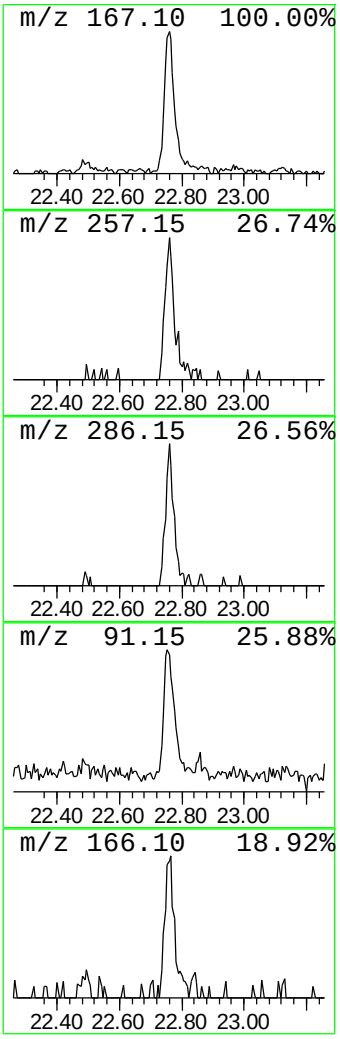
Peak Number 14 5-Methoxy-3,7-dihydroxyflav... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.76	2.41 ng	102969	Chrysene-d12	21.81

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	5-Methoxy-3,7-dihydroxyflavanone	286	C16H14O5	087620-04-0	50
2	5-Fluoro-2-methyl-benzothiazole	167	C8H6FNS	000399-75-7	49
3	2-Propanone, 1,1-diphenyl-	210	C15H14O	000781-35-1	49
4	Benzene, 1,1',1''-(1-ethanyl-2-y...	258	C20H18	001520-42-9	47
5	Acetamide, 2,2-diphenyl-N-(4-chl...	366	C20H15ClN2O3	1000295-48-3	47





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Instrument : BNA_G
ClientSampleId : MH-02

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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST11.L
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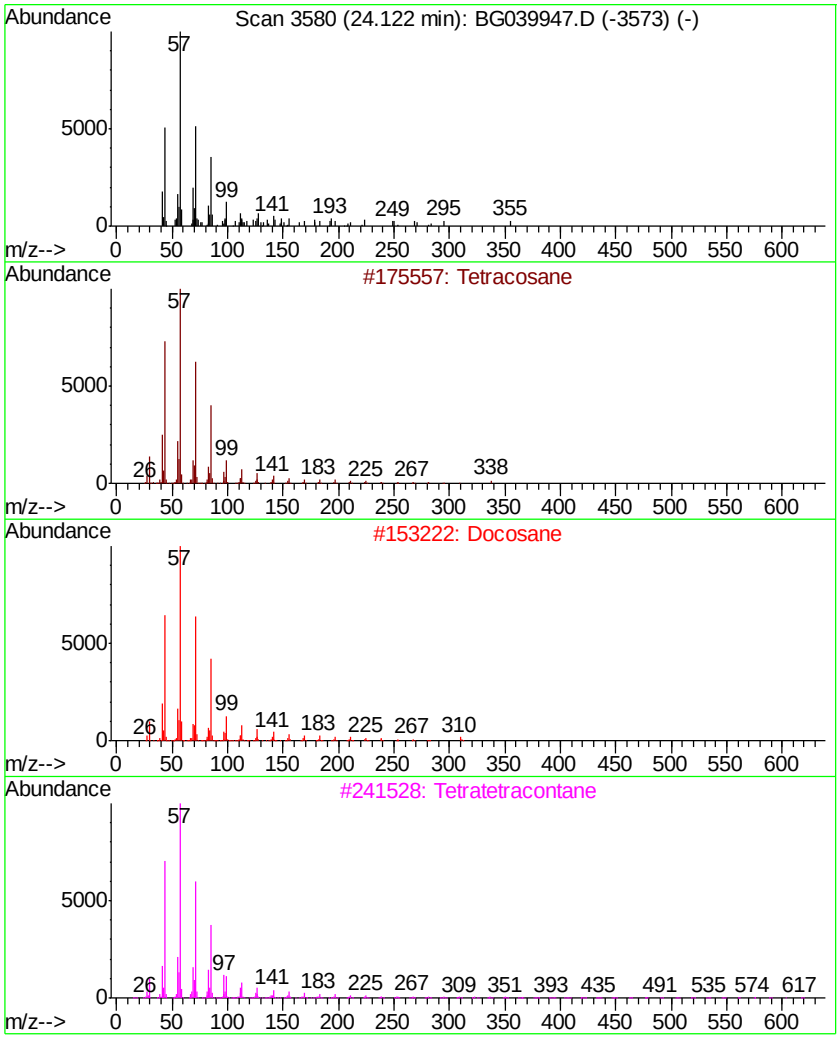
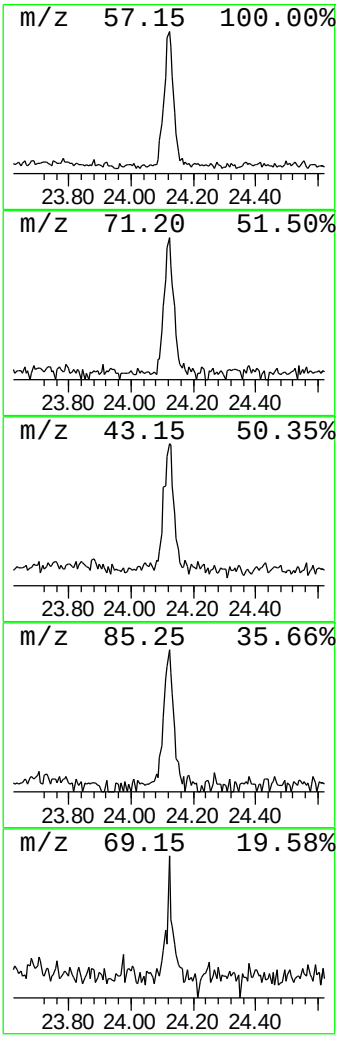
Peak Number 15 Tetracosane Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.12	2.09 ng	90510	Perylene-d12	25.17

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tetracosane	338	C24H50	000646-31-1	83
2	Docosane	310	C22H46	000629-97-0	83
3	Tetratetracontane	619	C44H90	007098-22-8	83
4	Hexacosane	366	C26H54	000630-01-3	83
5	Nonadecane	268	C19H40	000629-92-5	80





Tentatively Identified Compound (LSC) summary

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Data File : BG039947.D
Acq On : 15 Mar 2019 18:49
Operator : JU/SJ
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Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : MH-02

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG030419.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
Butane, 2-methoxy...	3.27	52.1	ng	723621	1	8.15	277582	20.0
2-Pentanone, 4-hy...	5.23	2.3	ng	31445	1	8.15	277582	20.0
unknown7.67	7.67	48.5	ng	673236	1	8.15	277582	20.0
Bicyclo[2.2.1]hep...	9.86	3.1	ng	68146	2	10.97	436660	20.0
(+)-Borneol	10.73	3.0	ng	66363	2	10.97	436660	20.0
Terpinen-4-ol	10.82	2.3	ng	51050	2	10.97	436660	20.0
7-Methoxymethyl-2...	15.28	18.5	ng	548357	3	14.78	592474	20.0
3,4,5-trimethoxyp...	18.26	3.8	ng	151345	4	17.52	789770	20.0
unknown18.52	18.52	19.5	ng	768908	4	17.52	789770	20.0
unknown18.73	18.73	3.2	ng	127638	4	17.52	789770	20.0
(R-(R*,R*))-4-(1,...	19.10	4.4	ng	172135	4	17.52	789770	20.0
5-Methoxy-3,7-dih...	22.76	2.4	ng	102969	5	21.81	855170	20.0
Tetracosane	24.12	2.1	ng	90510	6	25.17	867430	20.0