

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG062718\
 Data File : BG035463.D
 Acq On : 28 Jun 2018 2:38
 Operator : JU/SJ
 Sample : J3695-02
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 LIC-FIELDBLANK-1-06-20-18

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-BG060718.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	5.161	314	319	326	rBV	49954	77950	1.28%	0.348%
2	5.631	391	399	408	rBV	516911	857588	14.11%	3.834%
3	7.212	660	668	679	rBV	331021	572210	9.42%	2.558%
4	7.588	723	732	743	rBV	1009729	1738126	28.60%	7.771%
5	8.052	804	811	818	rBV2	177346	296834	4.88%	1.327%
6	8.375	858	866	878	rVB	803307	1393093	22.92%	6.228%
7	9.215	1000	1009	1017	rBV	808913	1429914	23.53%	6.393%
8	10.854	1280	1288	1297	rBV	218285	390202	6.42%	1.744%
9	13.297	1697	1704	1714	rBV	2118966	3239390	53.30%	14.482%
10	14.672	1930	1938	1948	rBV2	368426	602091	9.91%	2.692%
11	16.158	2184	2191	2209	rBV	1890648	3035675	49.95%	13.572%
12	17.415	2398	2405	2415	rBV	540971	812274	13.37%	3.631%
13	20.024	2842	2849	2858	rBV	4521599	6077096	100.00%	27.169%
14	21.680	3123	3131	3138	rBV	602164	969237	15.95%	4.333%
15	24.899	3669	3679	3692	rBV3	294971	876099	14.42%	3.917%

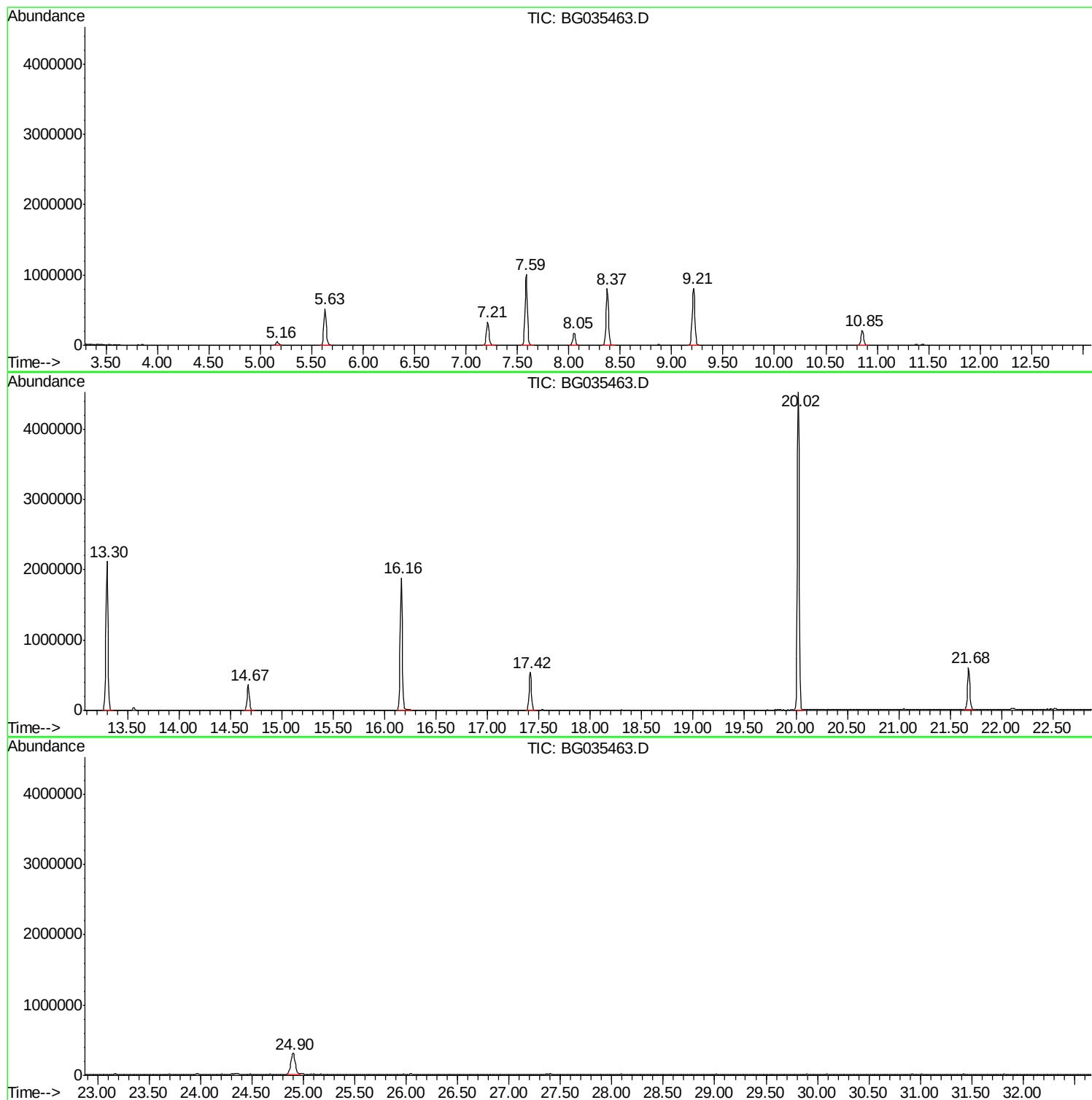
Sum of corrected areas: 22367779

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Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG060718.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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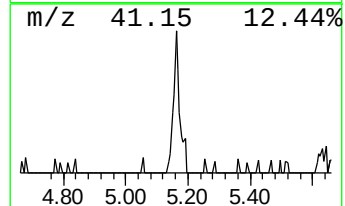
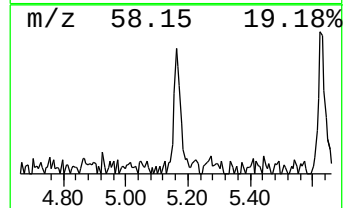
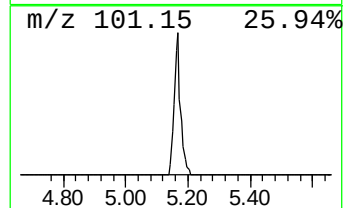
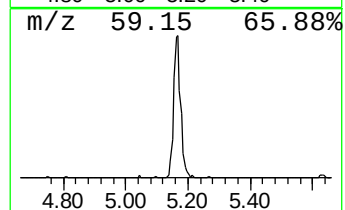
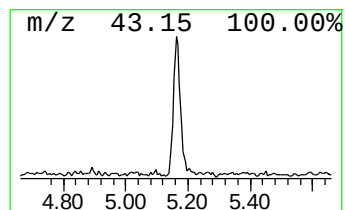
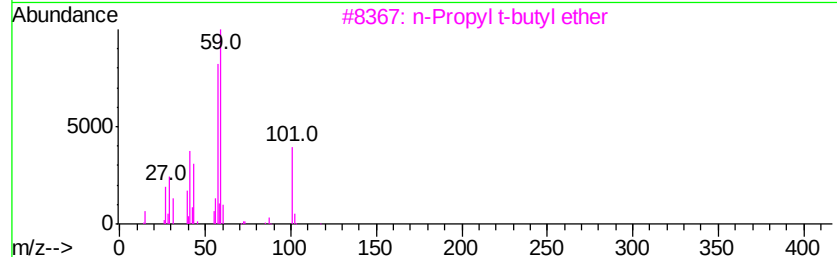
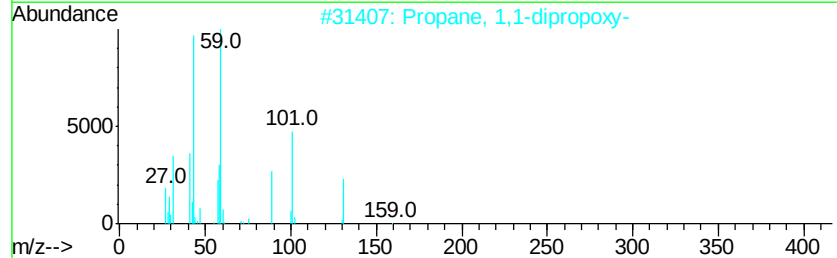
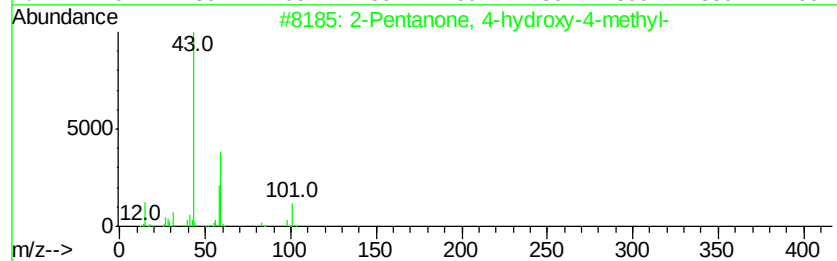
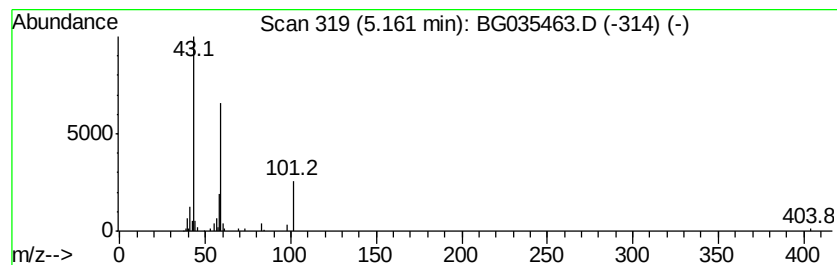
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.16	5.25 ng	77950	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	45
2			Propane, 1,1-dipropoxy-	160	C9H20O2	004744-11-0	39
3			n-Propyl t-butyl ether	116	C7H16O	1000334-81-9	25
4			4-Hydroxy-3-hexanone	116	C6H12O2	004984-85-4	25
5			Hydrazine, 1,1-bis(1-methylethyl)-	116	C6H16N2	000921-14-2	25



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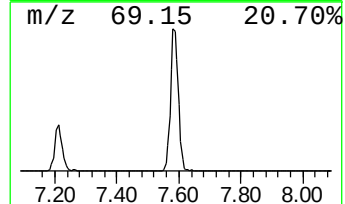
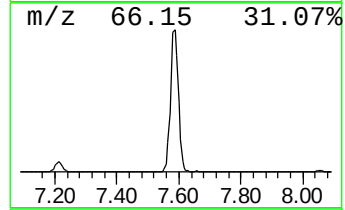
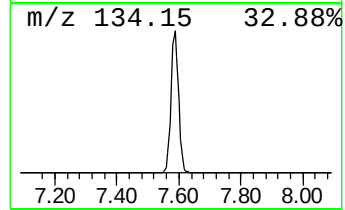
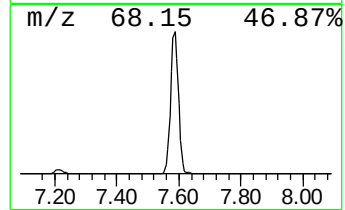
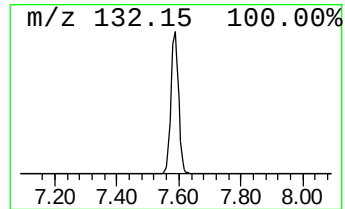
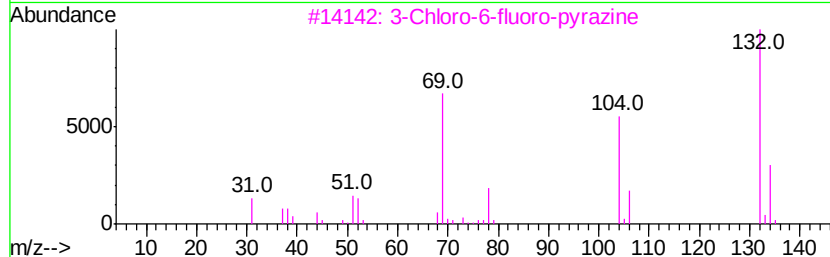
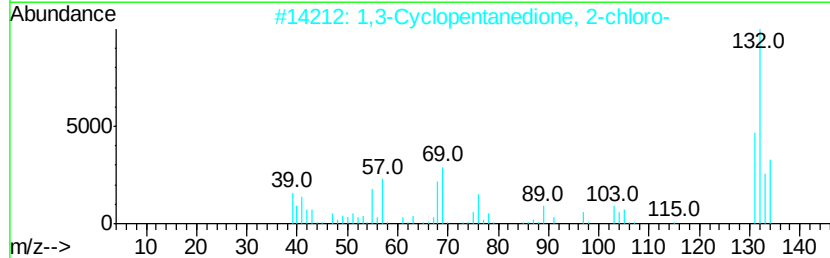
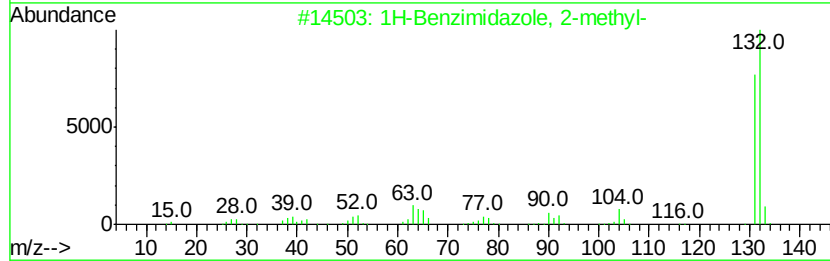
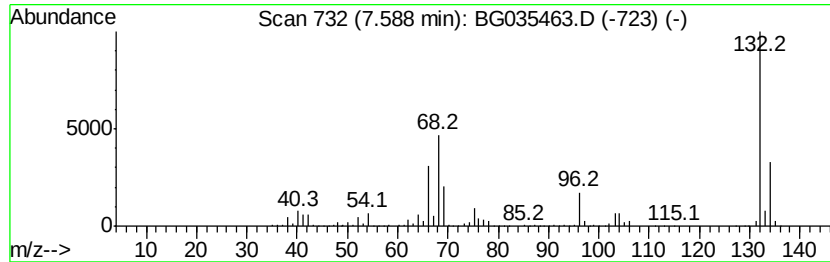
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Peak Number 2 unknown7.59 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.59	117.11 ng	1738130	1,4-Dichlorobenzene-d4	8.05

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	22
2		1,3-Cyclopentanedione, 2-chloro-	132	C5H5ClO2	014203-19-1	17
3		3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	16
4		5-Aminoindole	132	C8H8N2	005192-03-0	10
5		Cyclopropanamine, 1-phenyl-	133	C9H11N	1000351-26-2	10



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TIC Top Hit name	RT	EstConc	Units	Response	---Internal Standard---			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.16	5.3	ng	77950	1	8.05	296834	20.0
unknown7.59	7.59	117.1	ng	1738130	1	8.05	296834	20.0