

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG020719\
 Data File : BG039383.D
 Acq On : 7 Feb 2019 15:18
 Operator : JU/SJ
 Sample : PB116900BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB116900BL

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-BG012919.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	5.256	364	369	378	rVB	54062	83178	2.18%	0.437%
2	5.715	440	447	459	rVB	965219	1504839	39.45%	7.902%
3	7.312	711	719	732	rVB	983492	1666111	43.68%	8.749%
4	7.700	774	785	798	rBV	908186	1573916	41.27%	8.265%
5	8.176	859	866	874	rBV	187038	308065	8.08%	1.618%
6	8.499	913	921	930	rVB	651095	1150845	30.17%	6.043%
7	9.351	1058	1066	1078	rBV	550411	964197	25.28%	5.063%
8	11.002	1338	1347	1356	rBV	248107	452033	11.85%	2.374%
9	13.422	1752	1759	1767	rBV	1513187	2367189	62.06%	12.431%
10	14.802	1987	1994	2001	rBV2	394136	633730	16.62%	3.328%
11	16.283	2239	2246	2265	rBV	1100488	1654694	43.38%	8.689%
12	17.540	2454	2460	2468	rBV2	543208	826958	21.68%	4.343%
13	20.136	2896	2902	2908	rBV	3002968	3814108	100.00%	20.029%
14	21.834	3184	3191	3199	rBV2	624565	1047036	27.45%	5.498%
15	25.218	3755	3767	3779	rBV2	326749	996288	26.12%	5.232%

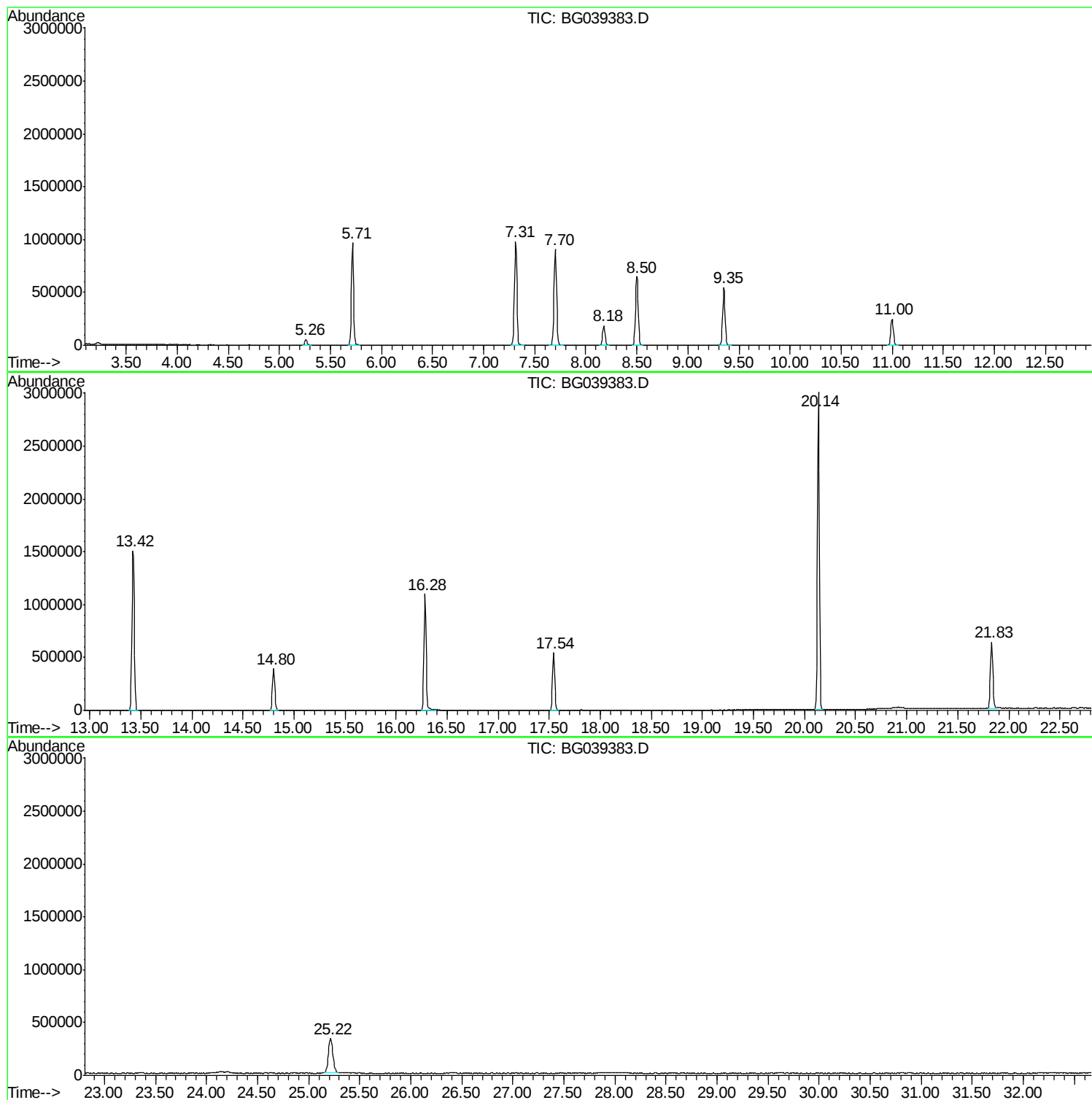
Sum of corrected areas: 19043187

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Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG012919.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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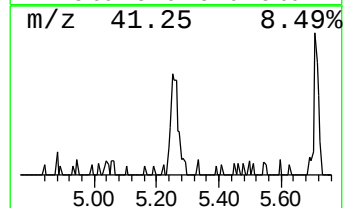
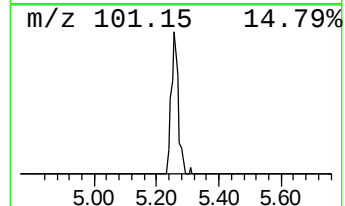
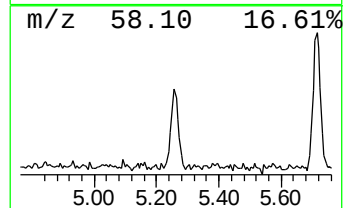
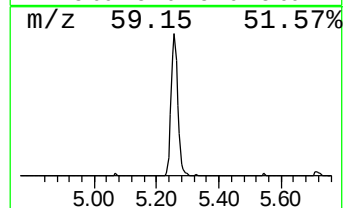
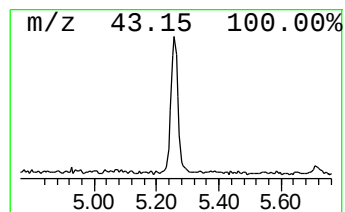
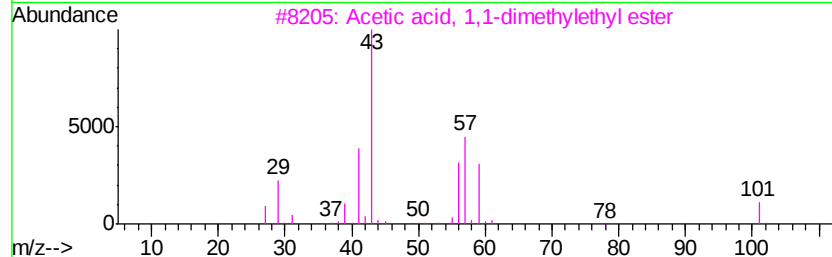
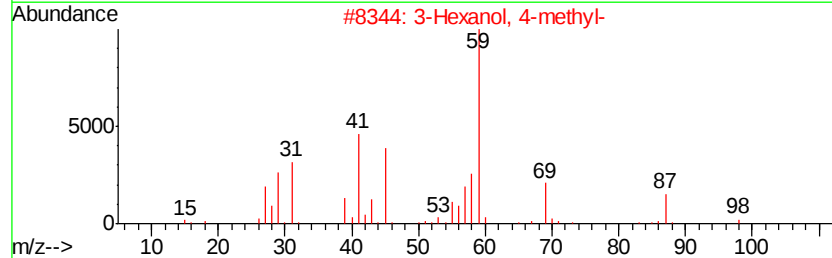
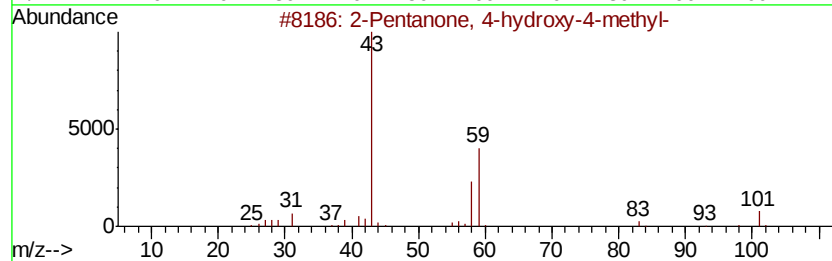
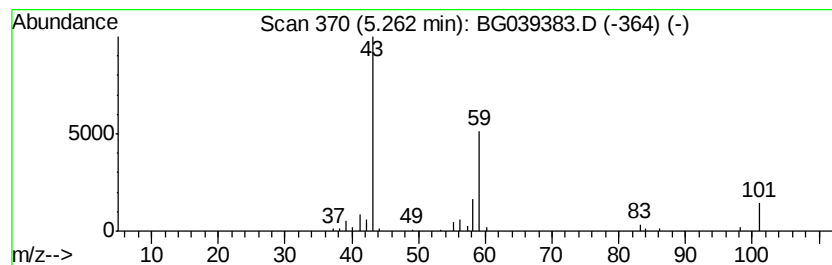
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.26	5.40 ng	83178	1,4-Dichlorobenzene-d4	8.18

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50
2			3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	28
3			Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	23
4			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	16
5			Guanidine	59	CH5N3	000113-00-8	9



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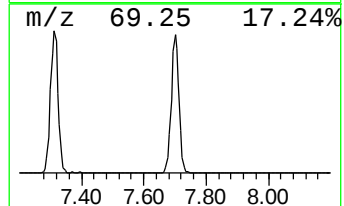
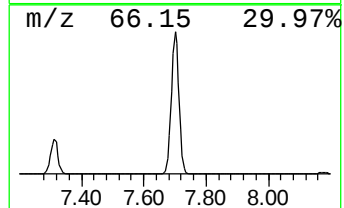
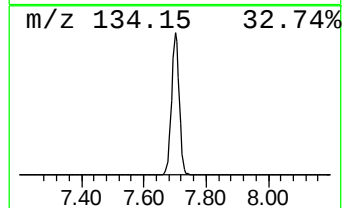
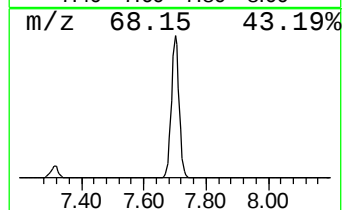
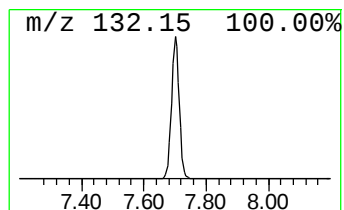
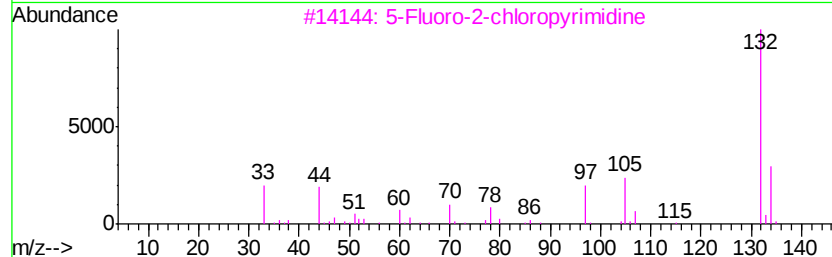
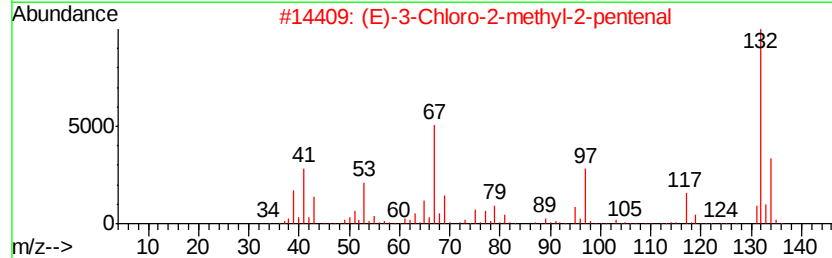
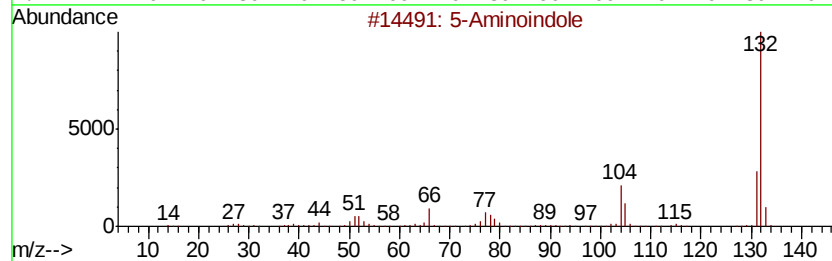
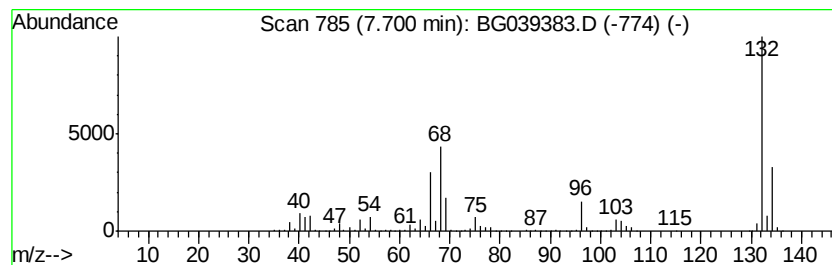
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Peak Number 2 unknown7.70 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.70	102.18 ng	1573920	1,4-Dichlorobenzene-d4	8.18

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	5-Aminoindole	132	C8H8N2	005192-03-0	22
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	16
3		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	12
4		(5-Methyl-2-pyridyl)acetonitrile	132	C8H8N2	1000241-93-9	10
5		1,3,5-Trifluorobenzene	132	C6H3F3	000372-38-3	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.26	5.4	ng	83178	1	8.18	308065	20.0
unknown7.70	7.70	102.2	ng	1573920	1	8.18	308065	20.0