

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG021119\
 Data File : BG039448.D
 Acq On : 12 Feb 2019 00:04
 Operator : JU/SJ
 Sample : PB116997BL
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB116997BL

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	5.251	362	368	379	rBV	63998	101666	2.06%	0.432%
2	5.709	439	446	459	rBV	1183995	1859518	37.66%	7.902%
3	7.307	710	718	730	rBV	1175160	2090818	42.35%	8.885%
4	7.695	776	784	797	rVB	1135784	1938416	39.26%	8.238%
5	8.171	858	865	873	rVB	180115	310144	6.28%	1.318%
6	8.494	912	920	928	rBV	794912	1347505	27.29%	5.726%
7	9.346	1057	1065	1074	rBV	663219	1203128	24.37%	5.113%
8	10.991	1338	1345	1353	rVB	274033	483714	9.80%	2.056%
9	13.417	1751	1758	1769	rBV	1968951	3127556	63.34%	13.291%
10	14.791	1986	1992	2004	rBV2	450405	764695	15.49%	3.250%
11	16.278	2238	2245	2265	rBV	1162734	1873350	37.94%	7.961%
12	17.535	2453	2459	2473	rBV2	623090	1003029	20.31%	4.263%
13	20.131	2895	2901	2907	rBV	3822691	4937534	100.00%	20.983%
14	21.829	3182	3190	3199	rBV2	681105	1176947	23.84%	5.002%
15	23.344	3442	3448	3456	rVB3	35950	67492	1.37%	0.287%
16	24.184	3585	3591	3598	rBV4	27658	64136	1.30%	0.273%
17	25.201	3752	3764	3780	rBV3	367282	1181502	23.93%	5.021%

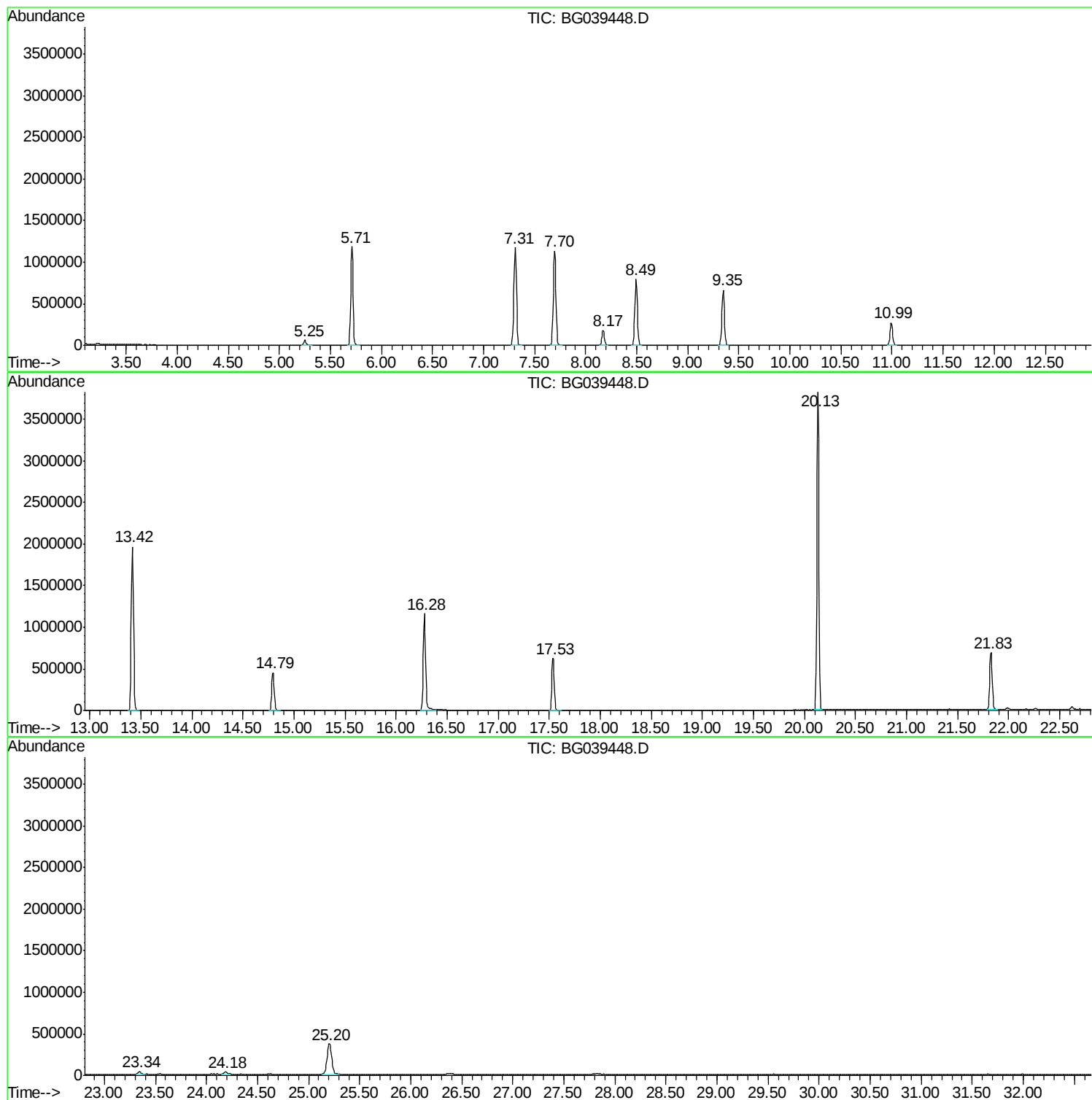
Sum of corrected areas: 23531150

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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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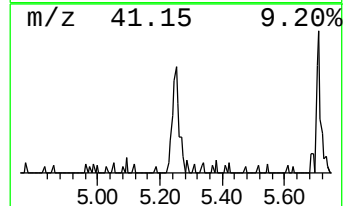
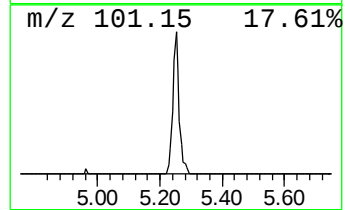
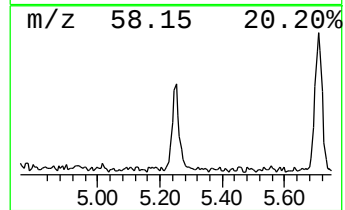
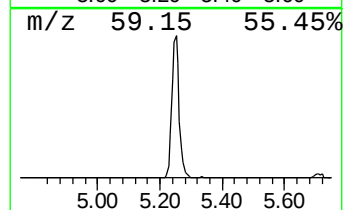
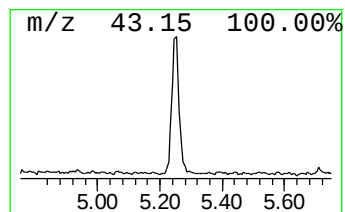
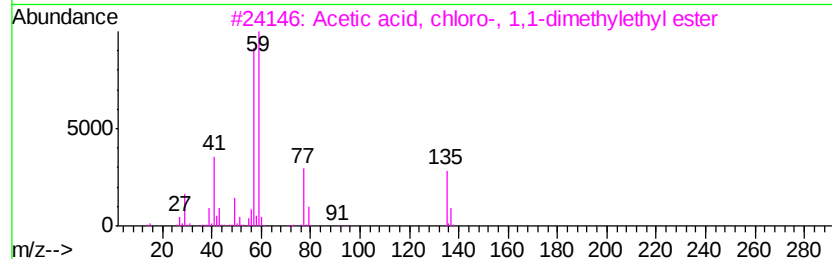
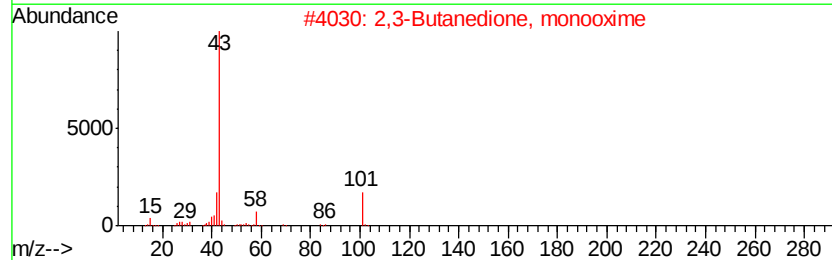
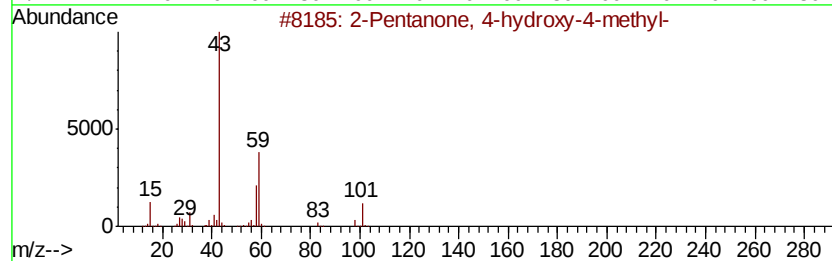
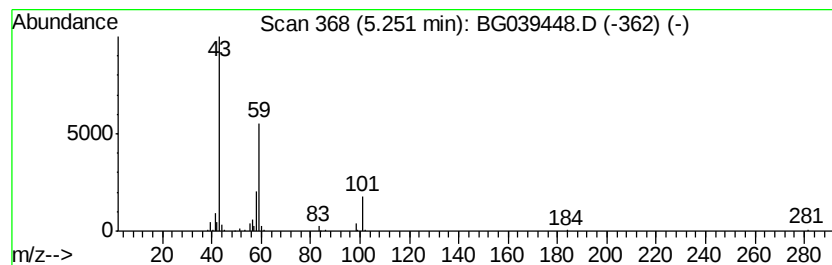
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.25	6.56 ng	101666	1,4-Dichlorobenzene-d4	8.17

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	64
2			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	12
3			Acetic acid, chloro-, 1,1-dimeth...	150	C6H11ClO2	000107-59-5	9
4			Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	9
5			Dimethadione	129	C5H7NO3	000695-53-4	9



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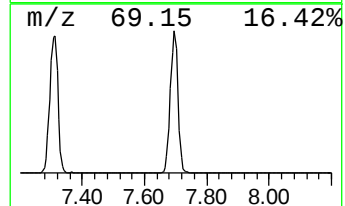
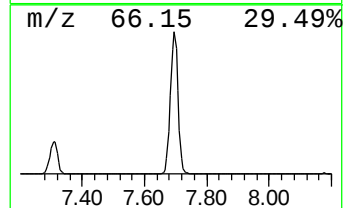
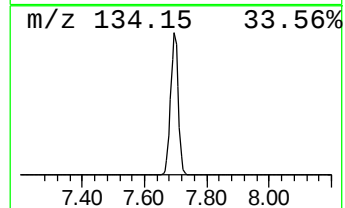
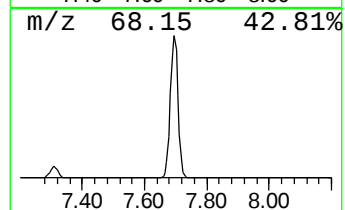
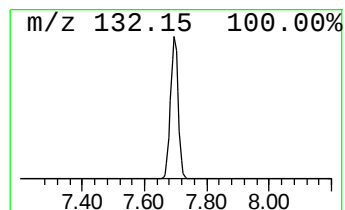
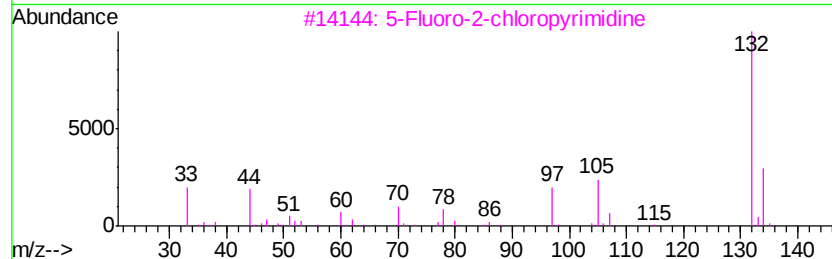
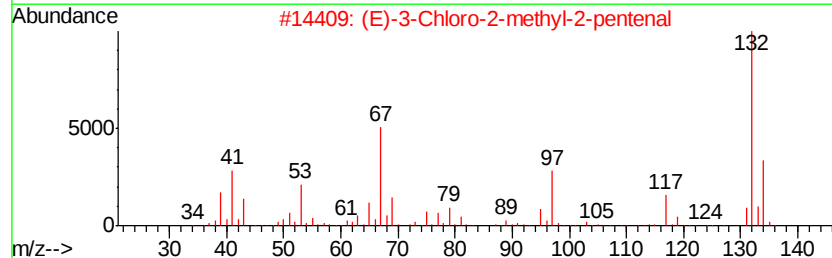
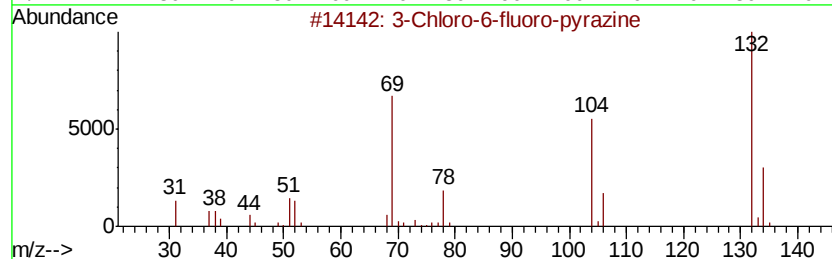
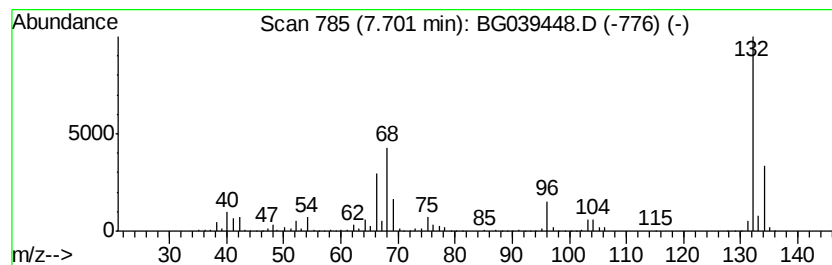
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Peak Number 2 unknown7.70 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.70	125.00 ng	1938420	1,4-Dichlorobenzene-d4	8.17

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	12
3		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	12
4		(5-Methyl-2-pyridyl)acetonitrile	132	C8H8N2	1000241-93-9	10
5		Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.25	6.6	ng	101666	1	8.17	310144	20.0
unknown7.70	7.70	125.0	ng	1938420	1	8.17	310144	20.0