

Data Path : Z:\HPCHEM1\BNA G\DATA\BG102317\
 Data File : BG029403.D
 Acq On : 23 Oct 2017 17:33
 Operator : SJ/JU
 Sample : PB103551BL
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB103551BL

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-BG101617.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	5.239	326	332	340	rBV	136652	202359	4.92%	1.051%
2	5.715	406	413	426	rBV	789845	1248060	30.35%	6.480%
3	7.319	677	686	701	rBV	760535	1336826	32.51%	6.940%
4	7.695	742	750	761	rBV	772364	1311166	31.89%	6.807%
5	8.165	822	830	837	rVB	187649	321472	7.82%	1.669%
6	8.488	878	885	893	rBV	574707	997662	24.26%	5.180%
7	9.328	1020	1028	1039	rBV	430750	766774	18.65%	3.981%
8	10.979	1302	1309	1319	rVB	262400	470692	11.45%	2.444%
9	13.406	1715	1722	1730	rBV	1312085	2111458	51.35%	10.962%
10	14.780	1950	1956	1965	rBV3	446359	722100	17.56%	3.749%
11	16.261	2201	2208	2228	rBV	1246116	1938116	47.13%	10.062%
12	17.518	2415	2422	2429	rBV2	708205	1073833	26.11%	5.575%
13	20.110	2857	2863	2873	rBV	3127898	4112040	100.00%	21.349%
14	21.790	3142	3149	3163	rBV2	777785	1339148	32.57%	6.953%
15	25.092	3700	3711	3727	rBV	440427	1309560	31.85%	6.799%

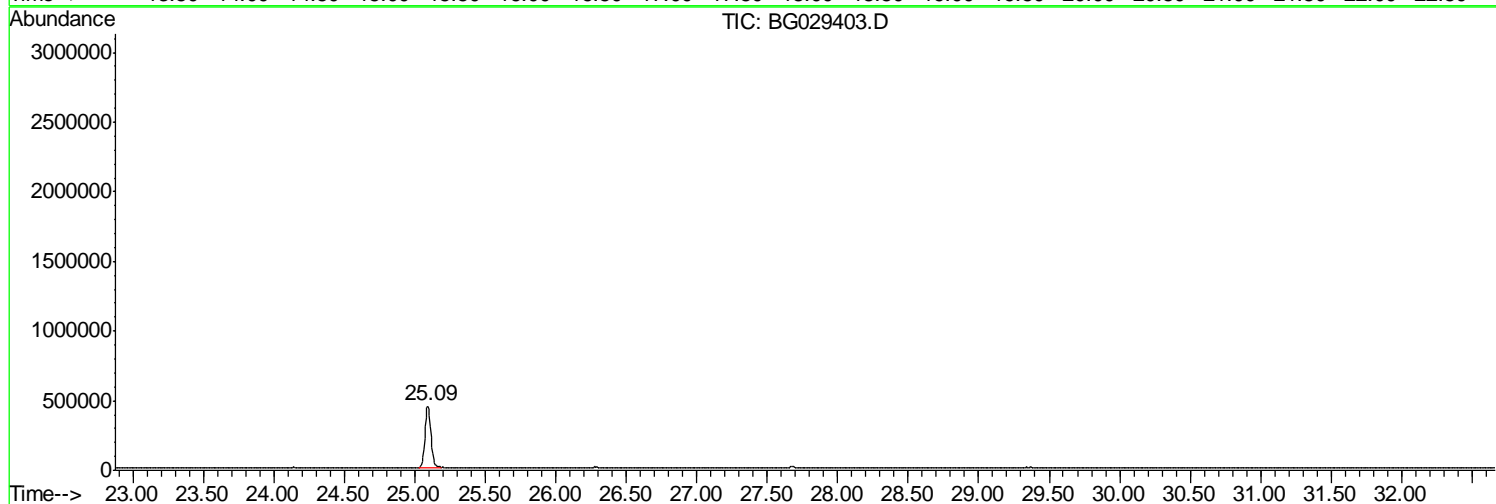
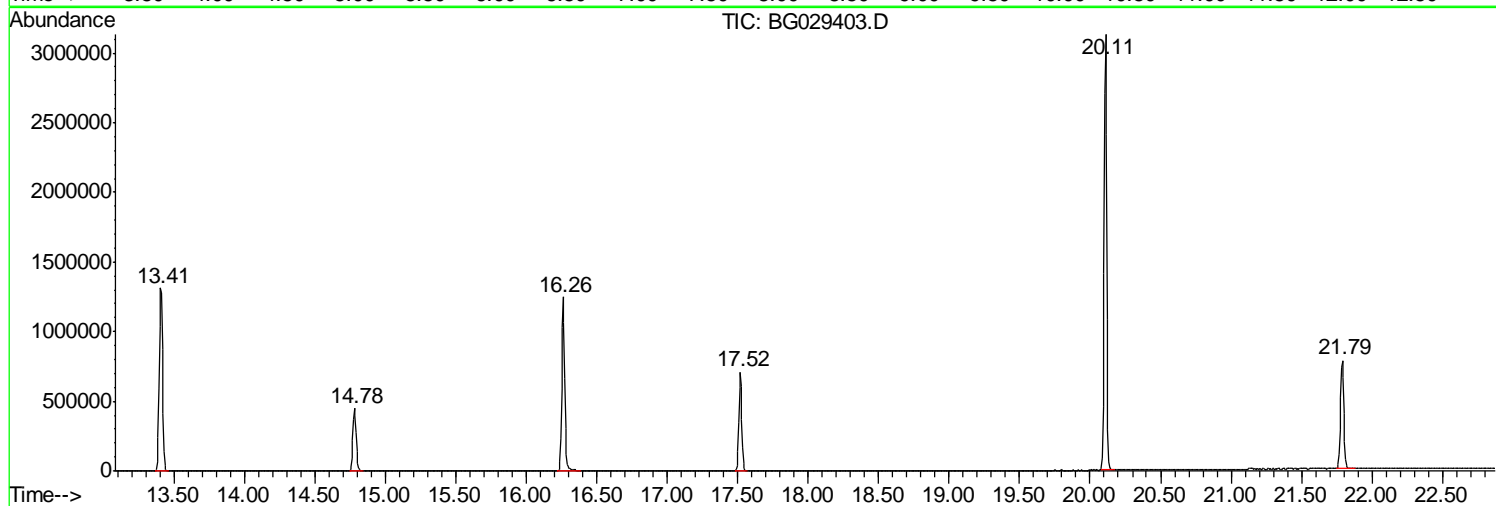
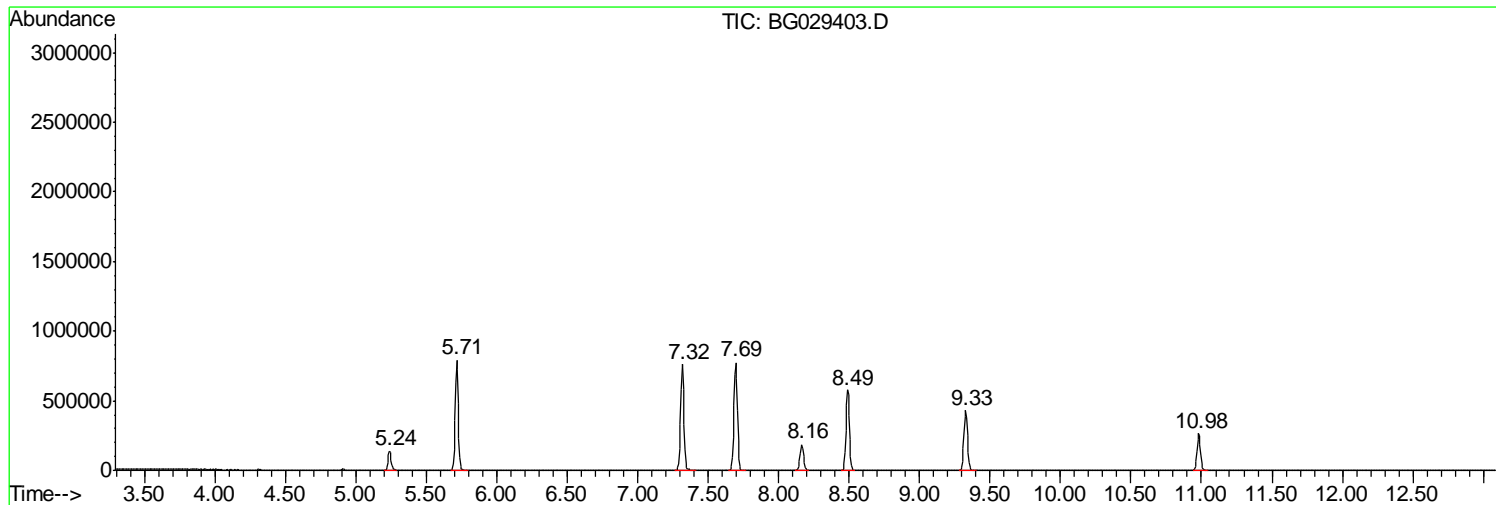
Sum of corrected areas: 19261266

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Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG101617.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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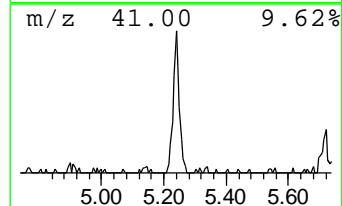
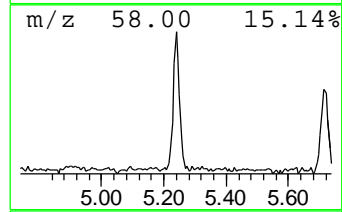
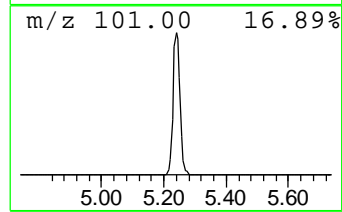
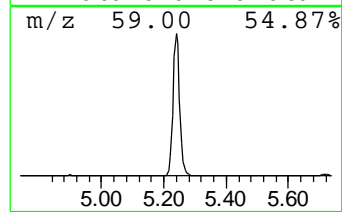
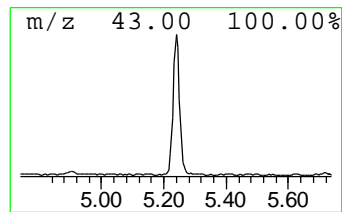
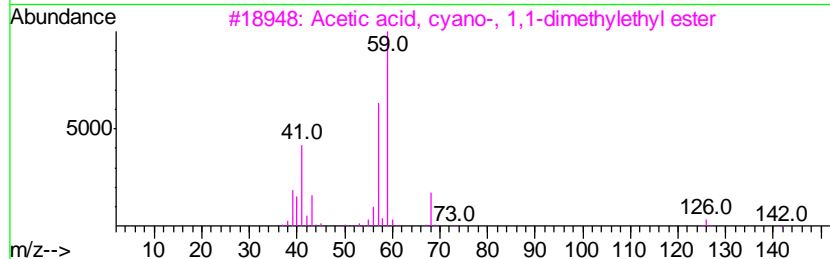
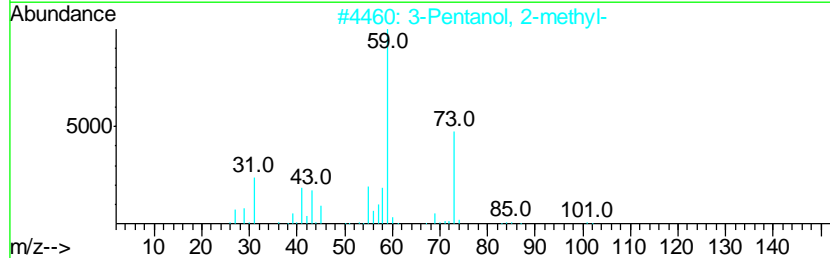
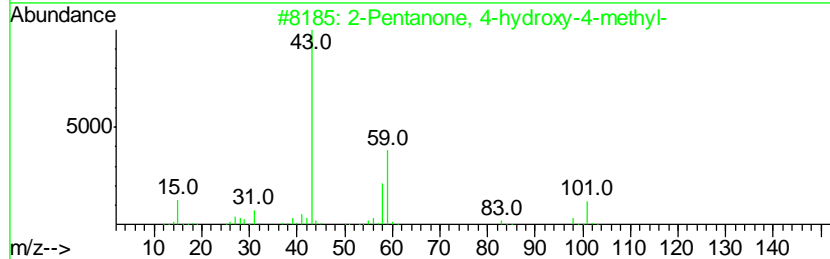
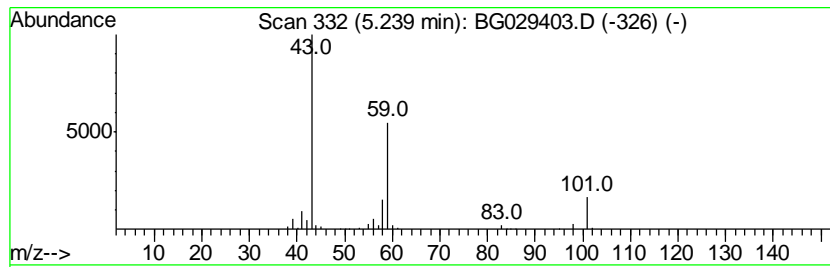
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 Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.24	12.59 ng	202359	1,4-Dichlorobenzene-d4	8.16

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	53
2		3-Pentanol, 2-methyl-	102	C6H14O	000565-67-3	40
3		Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	37
4		Propane, 2-methyl-2-(1-methyleth...	116	C7H16O	017348-59-3	36
5		2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	28



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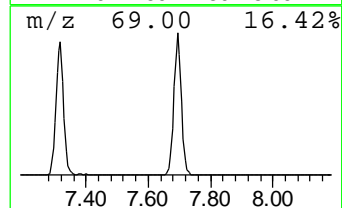
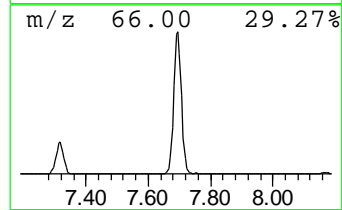
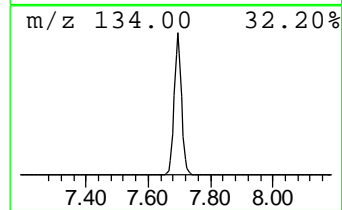
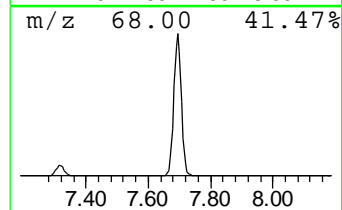
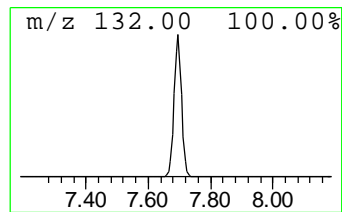
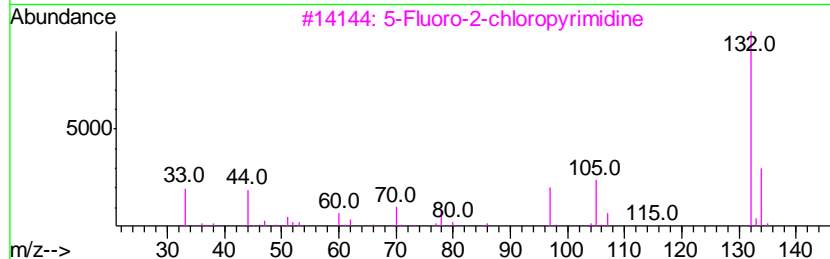
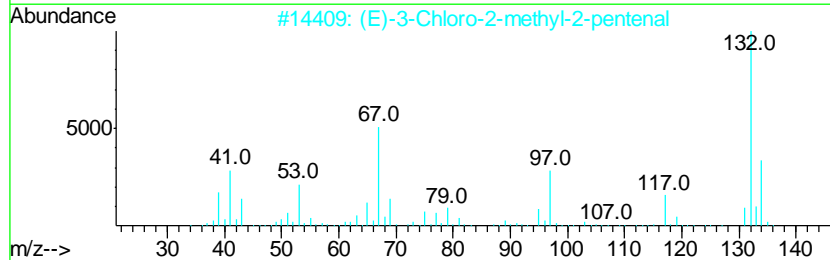
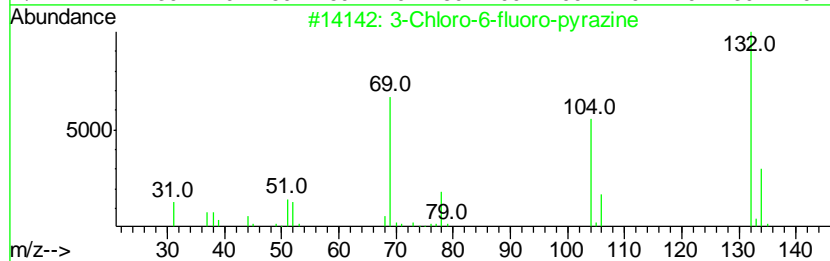
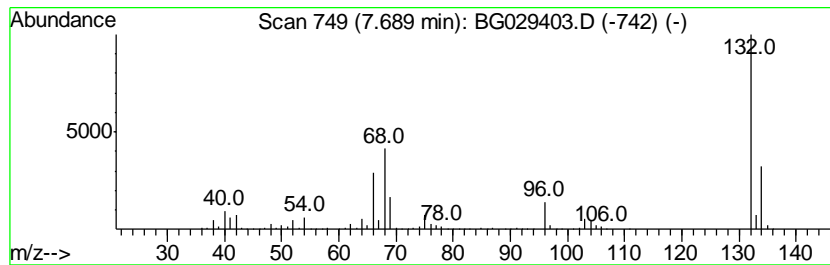
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 Peak Number 2 unknown7.69 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.69	81.57 ng	1311170	1,4-Dichlorobenzene-d4	8.16

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	27
3		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	12
4		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	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...	5.24	12.6	ng	202359	1	8.16	321472	20.0
unknown7.69	7.69	81.6	ng	1311170	1	8.16	321472	20.0