

Data Path : Z:\HPCHEM1\BNA G\DATA\BG102817\
 Data File : BG030531.D
 Acq On : 28 Oct 2017 17:22
 Operator : SJ/JU
 Sample : PB103686BL
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB103686BL

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	5.238	325	332	343	rVB	153257	235641	5.37%	1.108%
2	5.719	406	414	424	rBV	892661	1432572	32.64%	6.734%
3	7.318	678	686	698	rBV	872832	1496045	34.09%	7.032%
4	7.694	742	750	759	rBV	862562	1494348	34.05%	7.024%
5	8.164	823	830	838	rBV	226780	378438	8.62%	1.779%
6	8.493	878	886	896	rBV	670447	1168118	26.61%	5.491%
7	9.327	1019	1028	1036	rBV	489524	890730	20.29%	4.187%
8	10.984	1302	1310	1319	rBV	299182	542918	12.37%	2.552%
9	13.410	1715	1723	1738	rBV	1530684	2397590	54.63%	11.270%
10	14.785	1950	1957	1965	rBV2	489414	807852	18.41%	3.797%
11	16.266	2202	2209	2225	rBV	1263218	2072061	47.21%	9.740%
12	17.517	2416	2422	2432	rBV2	729307	1130016	25.75%	5.312%
13	20.108	2857	2863	2872	rBV	3412074	4389101	100.00%	20.631%
14	21.795	3142	3150	3160	rBV	847124	1440974	32.83%	6.773%
15	25.103	3702	3713	3728	rVB2	486341	1397566	31.84%	6.569%

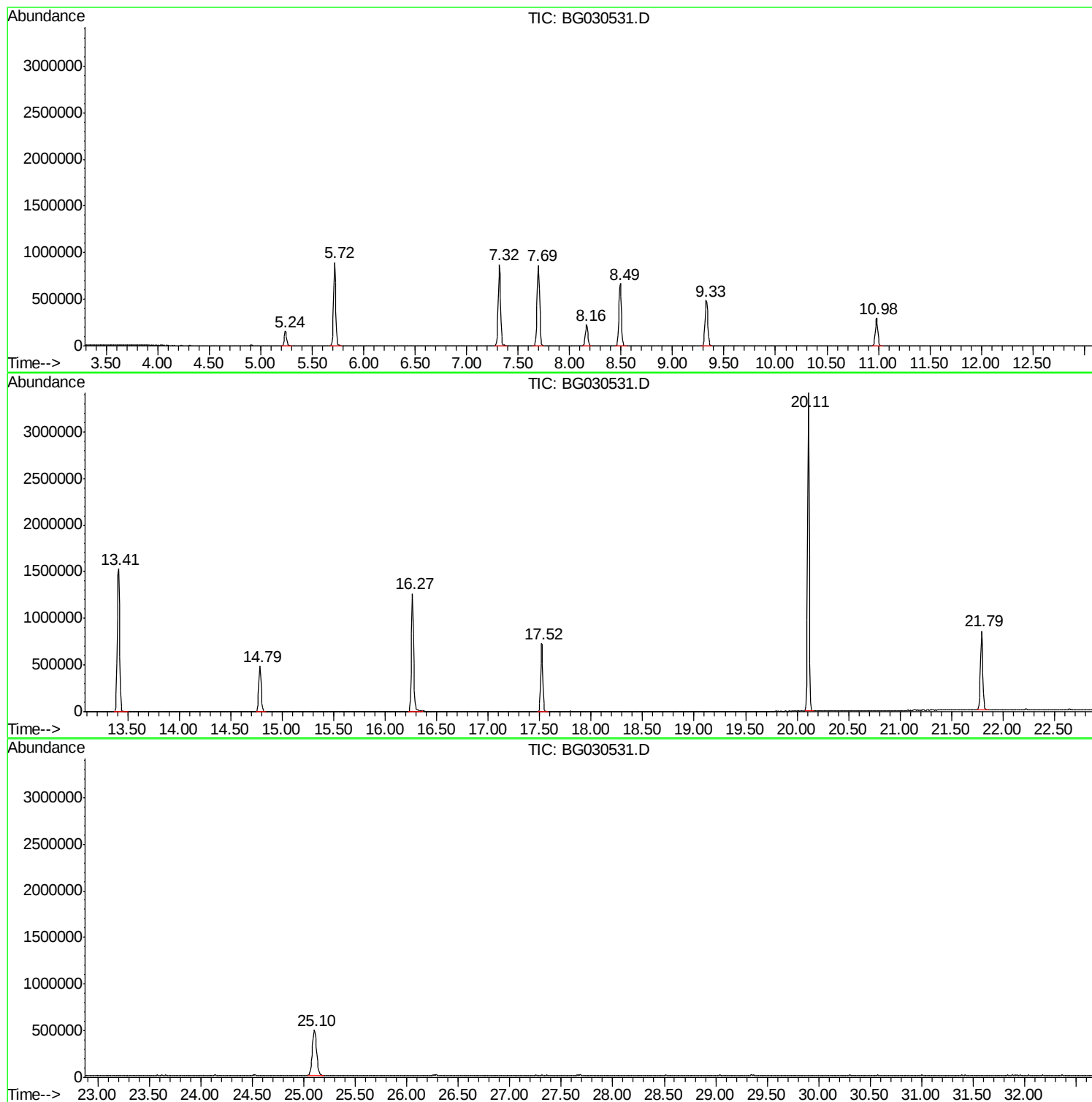
Sum of corrected areas: 21273970

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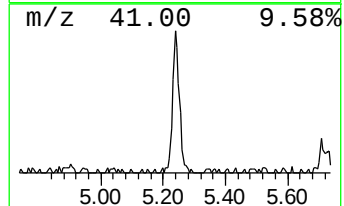
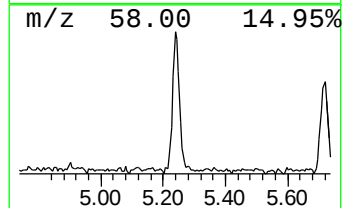
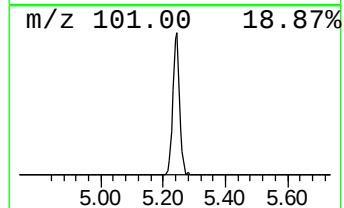
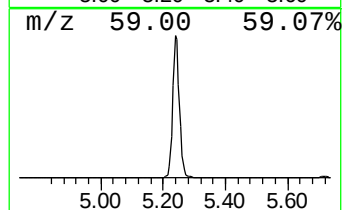
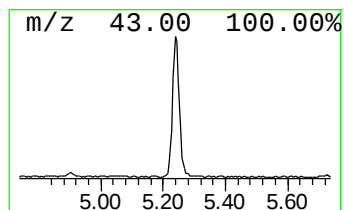
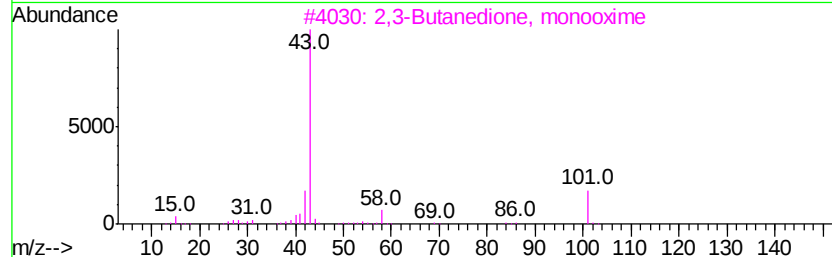
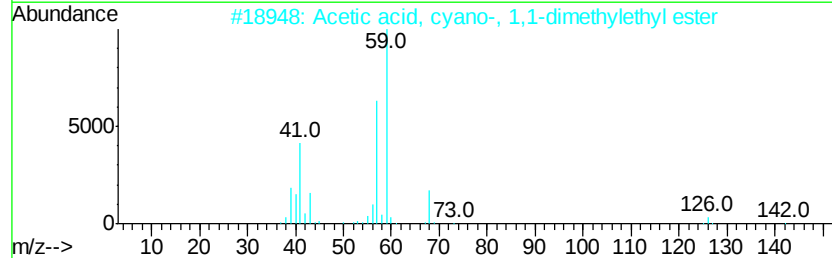
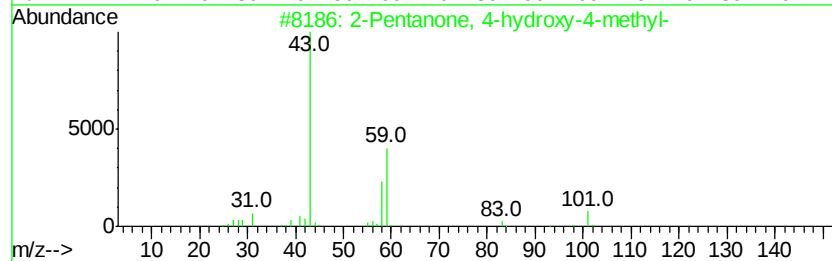
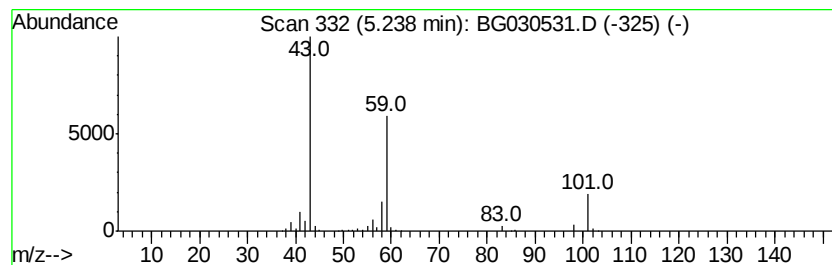
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.24	12.45 ng	235641	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	50
2		Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	23
3		2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	12
4		1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	10
5		Thiocyanic acid, propyl ester	101	C4H7NS	004251-16-5	9



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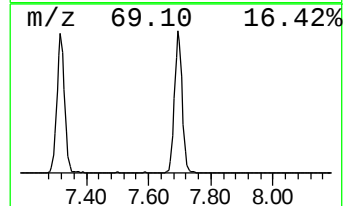
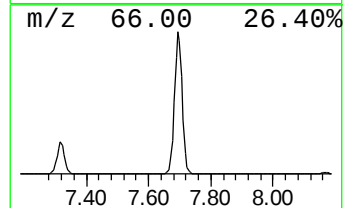
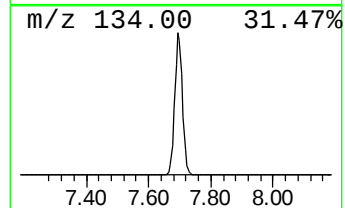
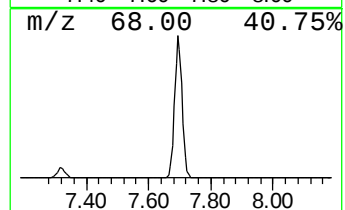
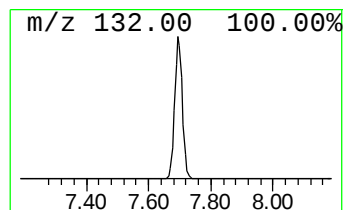
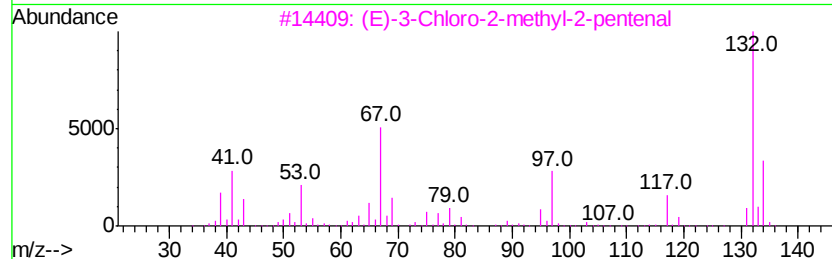
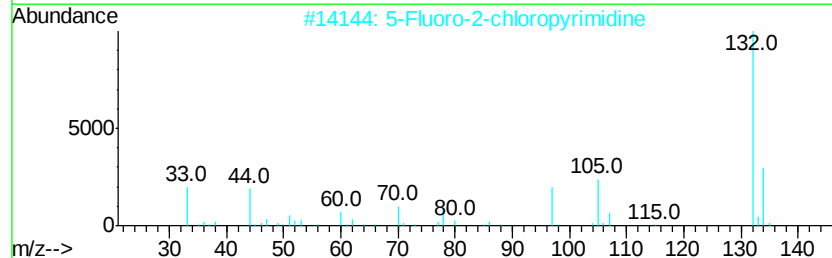
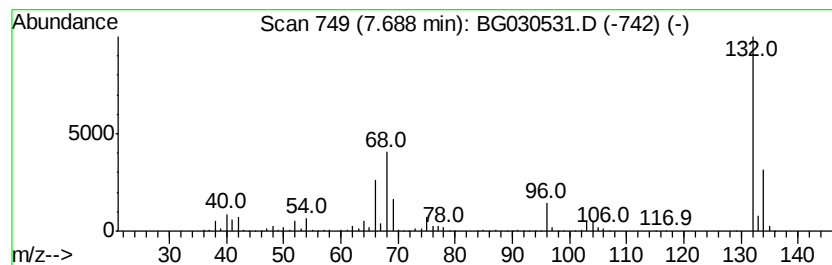
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Peak Number 2 unknown7.69 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.69	78.97 ng	1494350	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		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	16
3		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	12
4		5-Aminoindole	132	C8H8N2	005192-03-0	12
5		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.24	12.4	ng	235641	1	8.16	378438	20.0
unknown7.69	7.69	79.0	ng	1494350	1	8.16	378438	20.0