

Data Path : Z:\HPCHEM1\BNA G\DATA\BG011217\
 Data File : BG025484.D
 Acq On : 12 Jan 2017 16:14
 Operator : UM/SJ
 Sample : PB95939BL
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB95939BL

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-BG122116.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	5.260	405	412	422	rBV2	546817	904095	12.66%	2.911%
2	5.741	487	494	509	rBV	1268762	2214420	31.01%	7.130%
3	7.363	763	770	782	rBV	1237404	2347314	32.87%	7.558%
4	7.733	824	833	844	rBV	1260938	2261345	31.67%	7.281%
5	8.203	906	913	922	rBV	243915	421759	5.91%	1.358%
6	8.526	955	968	980	rBV	935794	1696763	23.76%	5.463%
7	9.384	1108	1114	1124	rBV	733018	1444053	20.22%	4.650%
8	11.035	1387	1395	1406	rBV	256847	561404	7.86%	1.808%
9	13.450	1799	1806	1818	rBV	2053552	3393722	47.52%	10.927%
10	14.831	2032	2041	2053	rBV2	443091	807631	11.31%	2.600%
11	16.317	2287	2294	2309	rBV	1946677	3348166	46.89%	10.780%
12	17.575	2502	2508	2520	rBV2	650302	1136731	15.92%	3.660%
13	20.154	2940	2947	2955	rBV	5221533	7141113	100.00%	22.993%
14	21.864	3230	3238	3249	rBV	882386	1701569	23.83%	5.479%
15	25.266	3807	3817	3831	rVB	562130	1677561	23.49%	5.401%

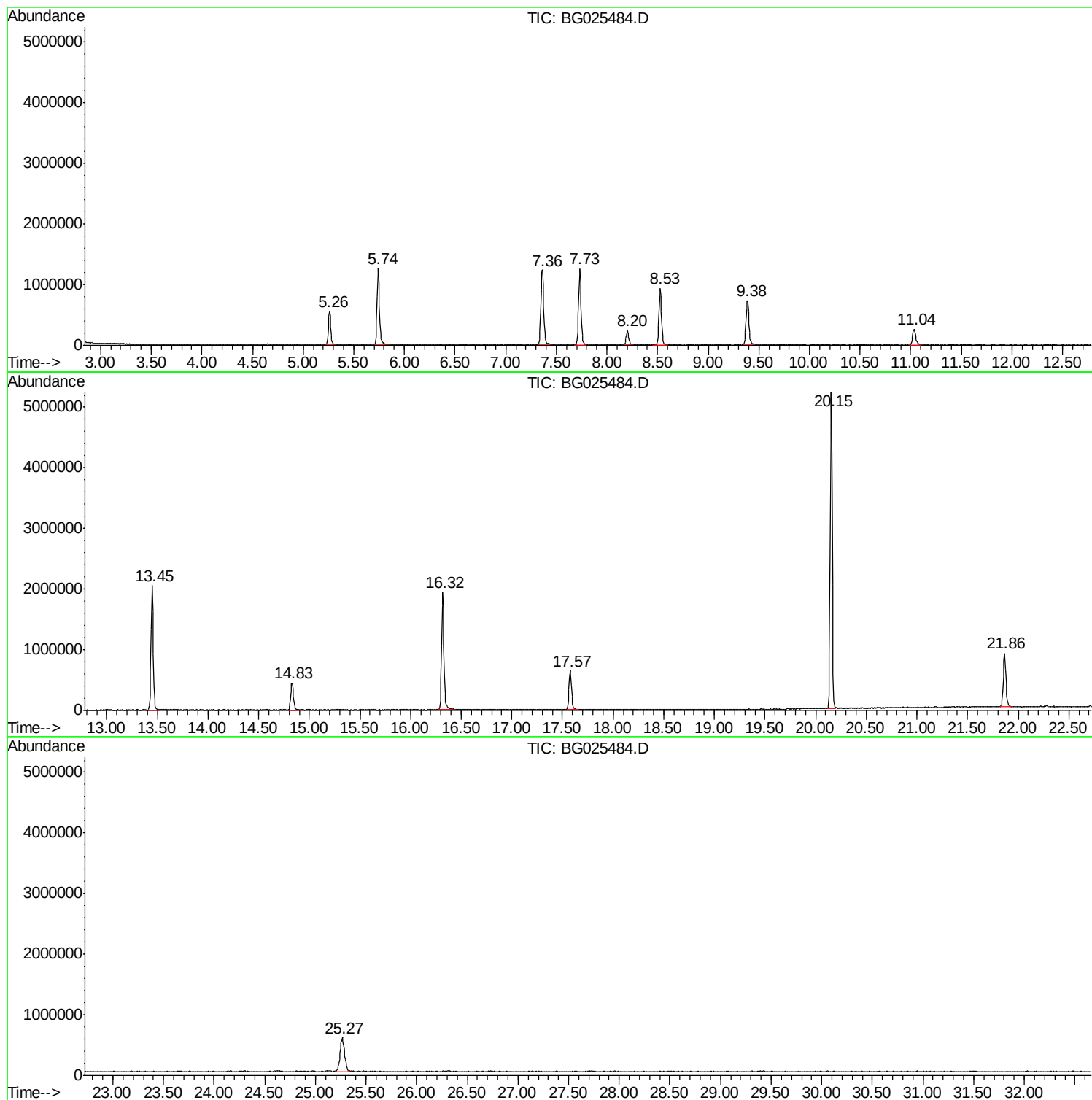
Sum of corrected areas: 31057646

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Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG122116.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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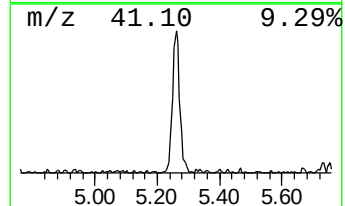
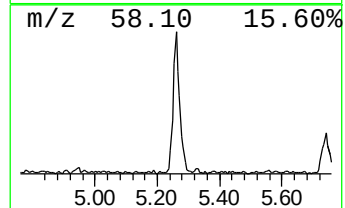
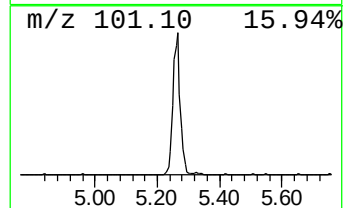
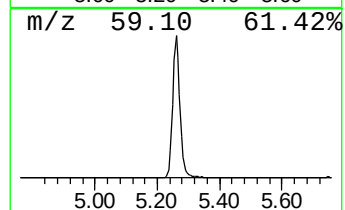
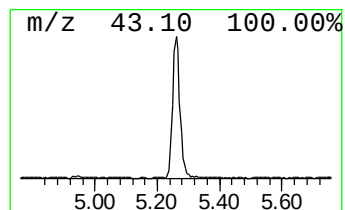
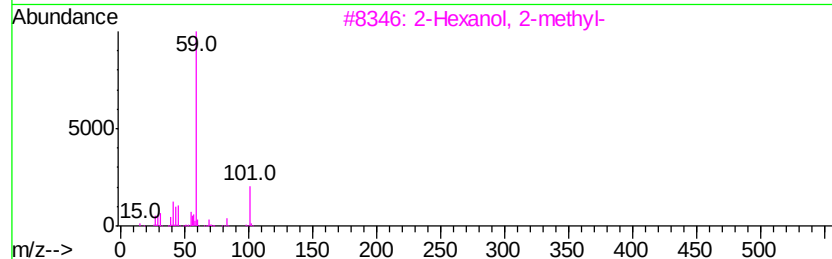
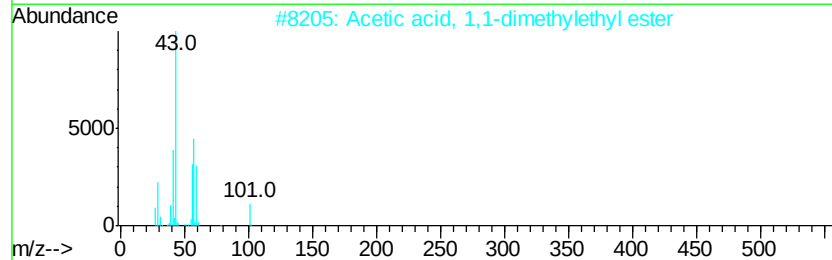
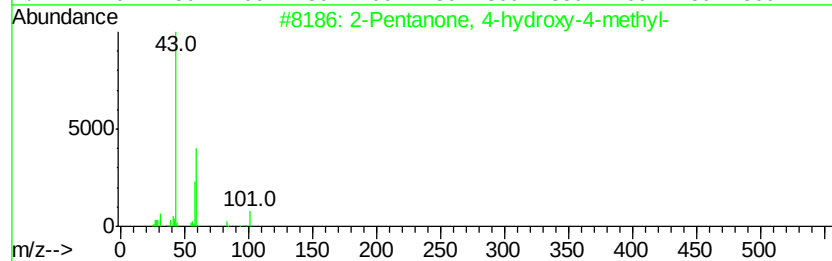
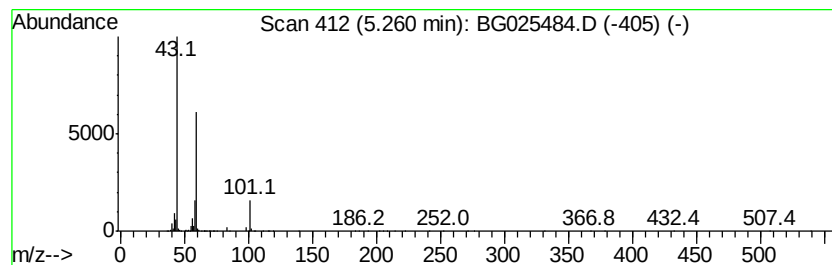
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.26	42.87 ng	904095	1,4-Dichlorobenzene-d4	8.20

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	64
2		Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	28
3		2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	28
4		Acetone	58	C3H6O	000067-64-1	9
5		1,3-Dioxane, 4,4-dimethyl-	116	C6H12O2	000766-15-4	9



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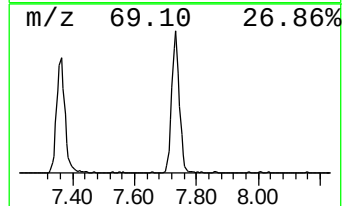
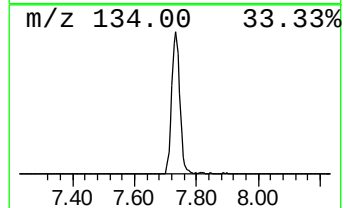
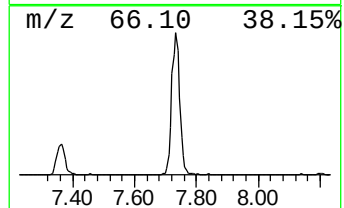
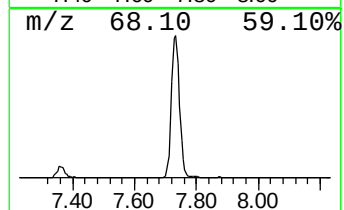
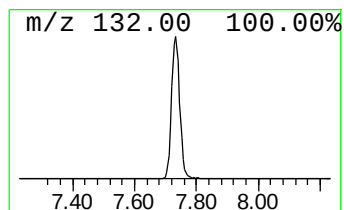
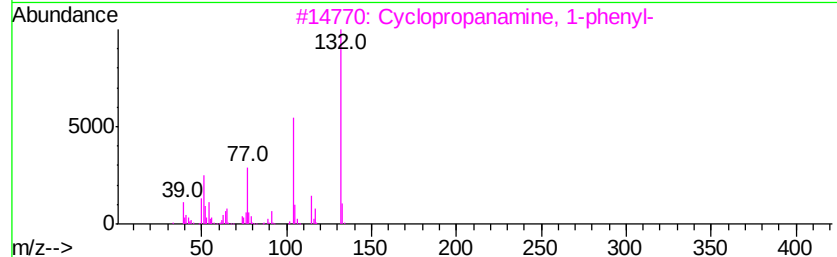
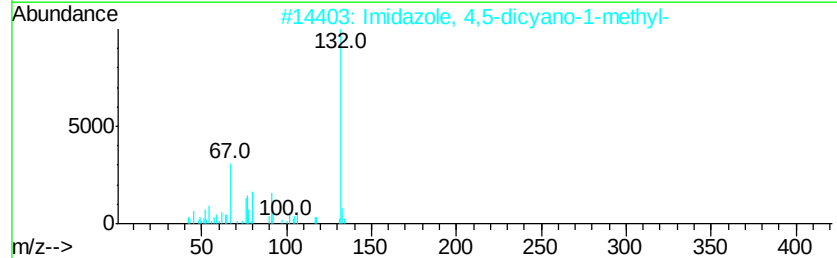
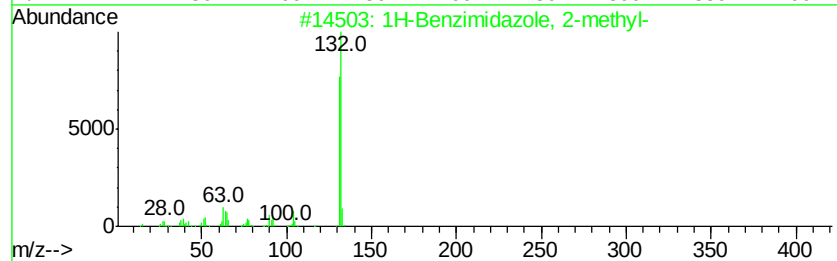
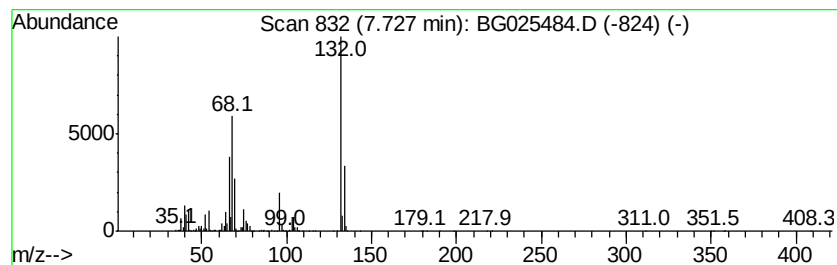
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Peak Number 2 unknown7.73 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.73	107.23 ng	2261350	1,4-Dichlorobenzene-d4	8.20

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	14
2		Imidazole, 4,5-dicyano-1-methyl-	132	C6H4N4	019485-35-9	11
3		Cyclopropanamine, 1-phenyl-	133	C9H11N	1000351-26-2	10
4		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	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	42.9	ng	904095	1	8.20	421759	20.0
unknown7.73	7.73	107.2	ng	2261350	1	8.20	421759	20.0