

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG062616\
 Data File : BG022630.D
 Acq On : 27 Jun 2016 4:44
 Operator : UM/SJ
 Sample : H3733-04
 Misc :
 ALS Vial : 82 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 MW-10

Integration Parameters: rteint.p
 Integrator: RTE
 Smoothing : ON
 Sampling : 1
 Start Thrs: 0.2
 Stop Thrs : 0
 Filtering: 5
 Min Area: 3 % of largest Peak
 Max Peaks: 100
 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
 Peak separation: 5

Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG062516.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	401	408	415	rBV	91262	149123	1.12%	0.257%
2	5.708	481	488	497	rBV	1209025	1947215	14.57%	3.358%
3	7.306	752	760	769	rBV	810599	1394857	10.44%	2.406%
4	7.688	816	825	835	rBV	2218182	3875239	29.00%	6.684%
5	8.158	899	905	914	rVB	539580	946880	7.09%	1.633%
6	8.487	953	961	968	rBV	1992276	3469948	25.97%	5.985%
7	9.333	1097	1105	1114	rBV	1863644	3322975	24.87%	5.731%
8	10.984	1379	1386	1394	rBV	732479	1370154	10.25%	2.363%
9	13.423	1793	1801	1807	rBV	4692478	7656119	57.29%	13.204%
10	14.798	2029	2035	2044	rVB2	1241709	2083857	15.59%	3.594%
11	16.290	2280	2289	2298	rBV	5200828	8190592	61.29%	14.126%
12	17.553	2497	2504	2513	rBV2	1823880	2892681	21.65%	4.989%
13	18.423	2648	2652	2658	rBV3	147800	227230	1.70%	0.392%
14	19.692	2863	2868	2874	rBV	194846	281273	2.10%	0.485%
15	19.798	2882	2886	2890	rBV	106578	154821	1.16%	0.267%
16	20.156	2941	2947	2953	rBV	9647027	13363878	100.00%	23.048%
17	21.848	3229	3235	3244	rVB	2041860	3415318	25.56%	5.890%
18	25.215	3799	3808	3824	rVB2	1116324	3239526	24.24%	5.587%

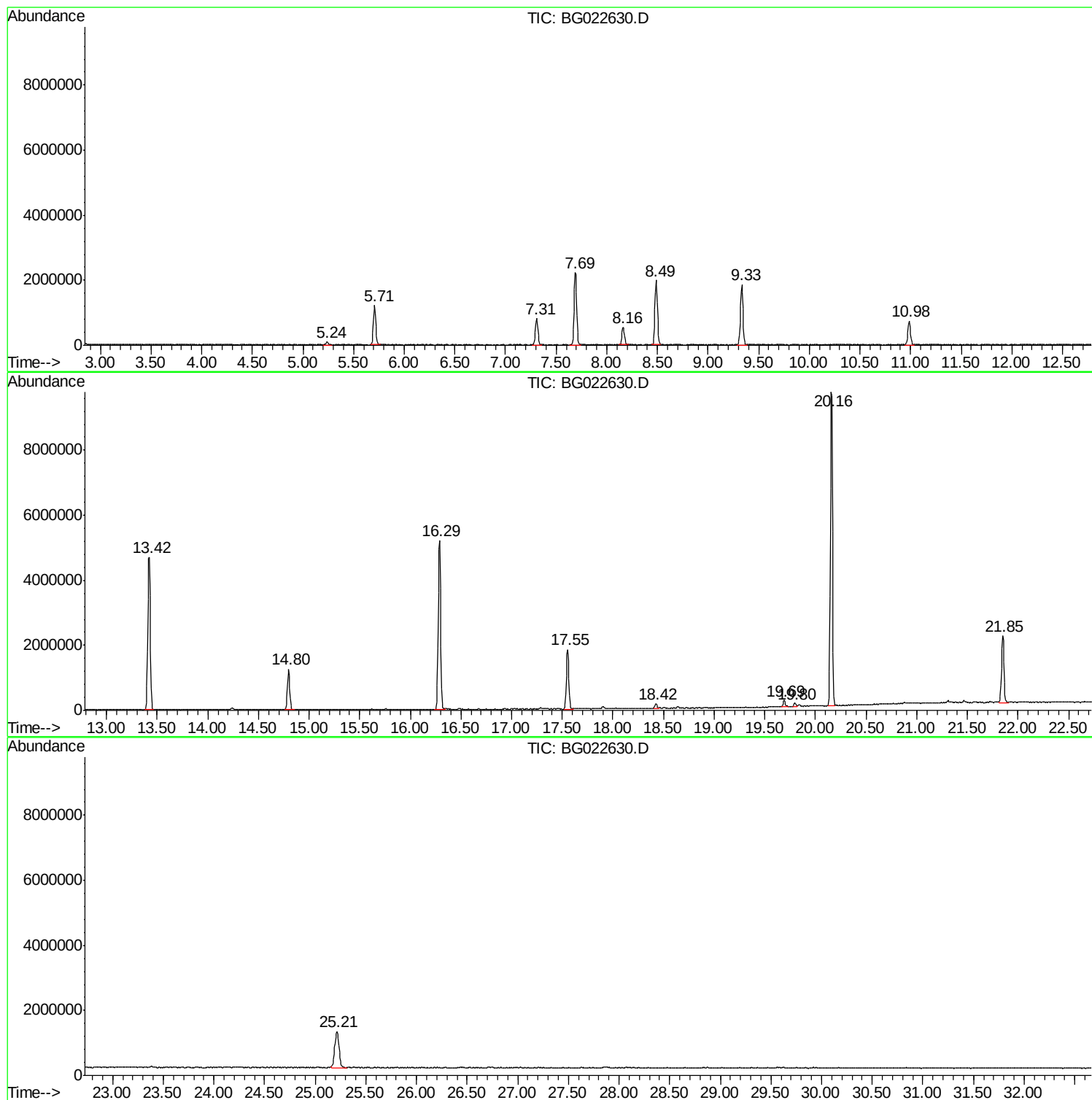
Sum of corrected areas: 57981686

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Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG062516.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P



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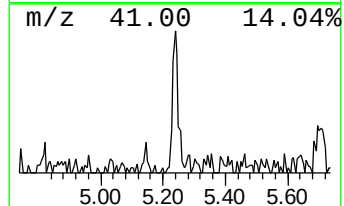
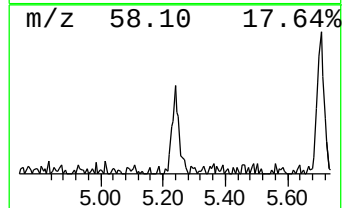
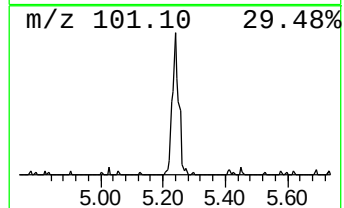
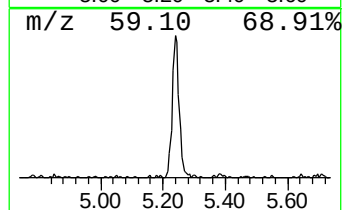
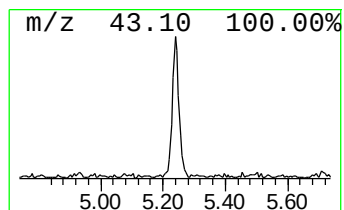
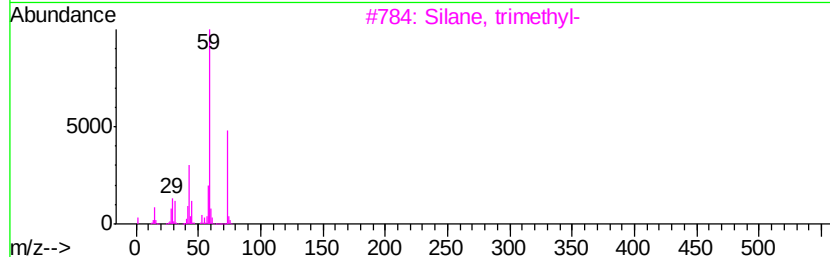
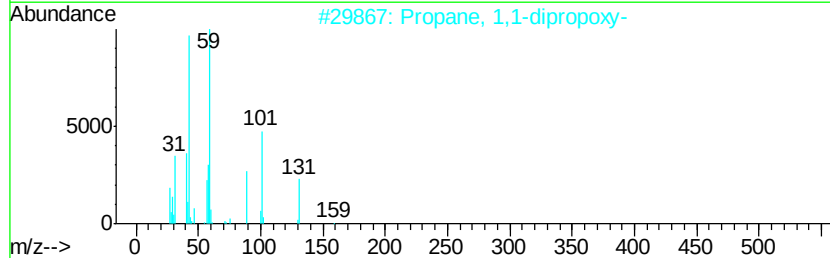
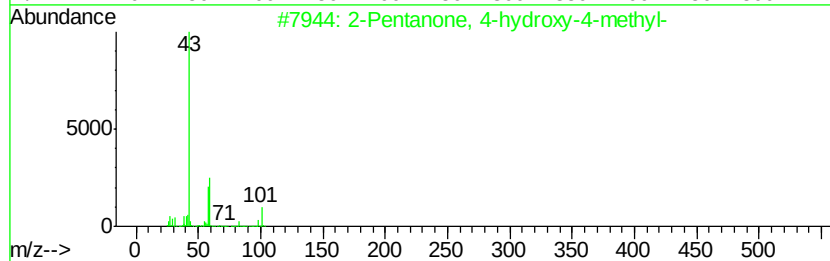
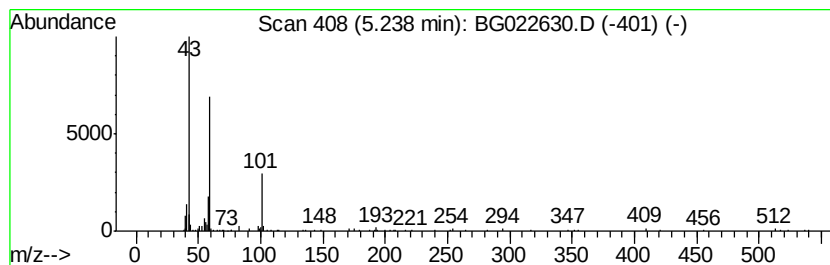
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.24	3.15 ng	149123	1,4-Dichlorobenzene-d4	8.16

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	59
2			Propane, 1,1-dipropoxy-	160	C9H20O2	004744-11-0	38
3			Silane, trimethyl-	74	C3H10Si	000993-07-7	23
4			2-Nonanone, 9-hydroxy-	158	C9H18O2	025368-56-3	23
5			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9



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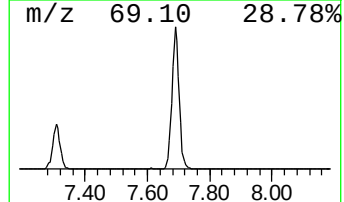
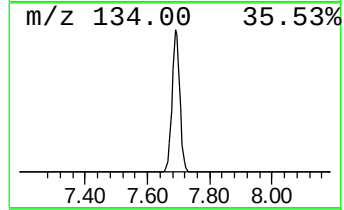
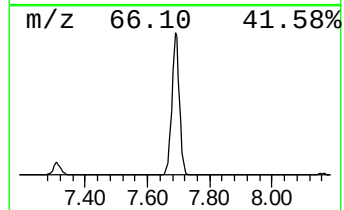
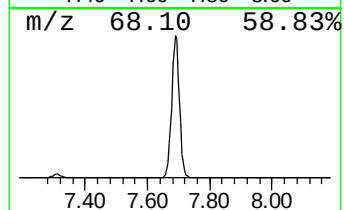
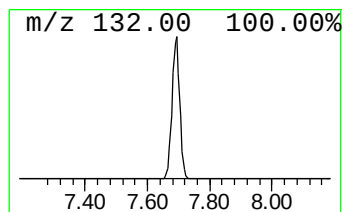
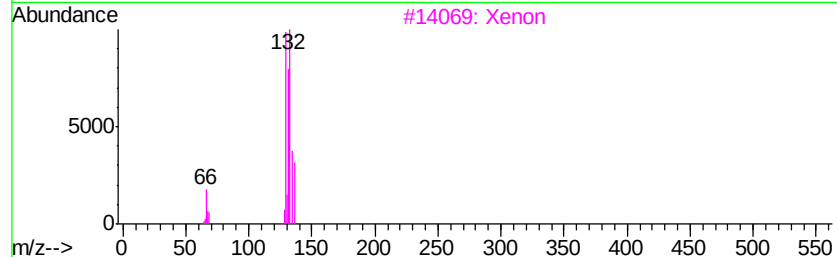
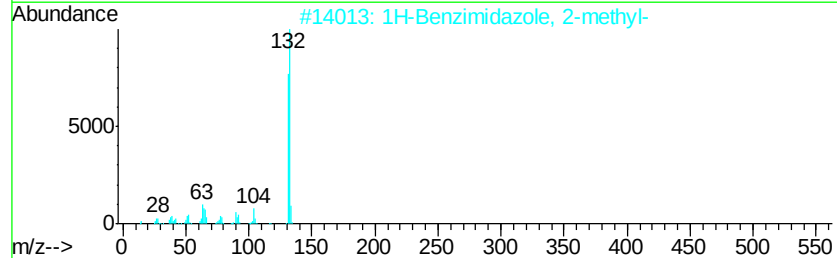
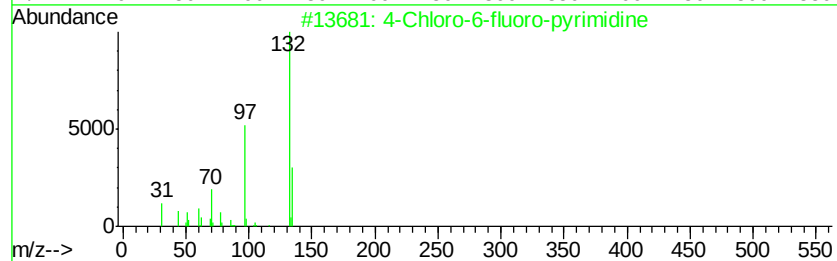
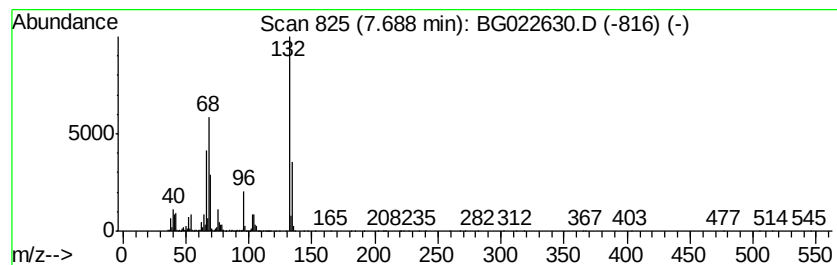
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Peak Number 2 unknown7.69 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.69	81.85 ng	3875240	1,4-Dichlorobenzene-d4	8.16

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	4-Chloro-6-fluoro-pyrimidine	132	C4H2ClFN2	051422-01-6	14
2		1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	10
3		Xenon	132	Xe	007440-63-3	9
4		Benzene, 1,3,5-trifluoro-	132	C6H3F3	000372-38-3	9
5		2H-Cyclopenta[d]pyridazine, 2-me...	132	C8H8N2	022291-85-6	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.24	3.1	ng	149123	1	8.16	946880	20.0
unknown7.69	7.69	81.8	ng	3875240	1	8.16	946880	20.0