

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA G\DATA\BG120817\
 Data File : BG031443.D
 Acq On : 9 Dec 2017 6:37
 Operator : SJ/JU
 Sample : I6744-07
 Misc :
 ALS Vial : 28 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PROS-2-E(4-5)

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-BG111017.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	5.204	321	326	332	rVB	37563	58777	1.50%	0.369%
2	5.686	402	408	425	rVB	516011	934428	23.85%	5.869%
3	7.290	674	681	701	rBV	573652	1113856	28.43%	6.996%
4	7.660	737	744	758	rBV	568575	1059329	27.04%	6.654%
5	8.124	817	823	834	rBV	127803	226614	5.78%	1.423%
6	8.447	871	878	888	rBV	435755	794934	20.29%	4.993%
7	9.293	1014	1022	1039	rBV	302727	641284	16.37%	4.028%
8	10.944	1295	1303	1312	rBV	147112	322376	8.23%	2.025%
9	13.371	1709	1716	1729	rBV	1110787	1910158	48.75%	11.998%
10	14.205	1851	1858	1873	rBV	74869	168339	4.30%	1.057%
11	14.752	1944	1951	1966	rVB2	323531	576091	14.70%	3.619%
12	15.568	2086	2090	2095	rVB2	32727	42162	1.08%	0.265%
13	16.238	2198	2204	2220	rBV	532247	1038963	26.52%	6.526%
14	16.426	2232	2236	2244	rVB	31082	47374	1.21%	0.298%
15	17.490	2411	2417	2431	rBV2	541168	911039	23.25%	5.722%
16	20.081	2852	2858	2868	rBV	2895975	3918267	100.00%	24.612%
17	21.767	3138	3145	3158	rBV2	604085	1117104	28.51%	7.017%
18	25.063	3694	3706	3730	rBV2	325082	1039353	26.53%	6.528%

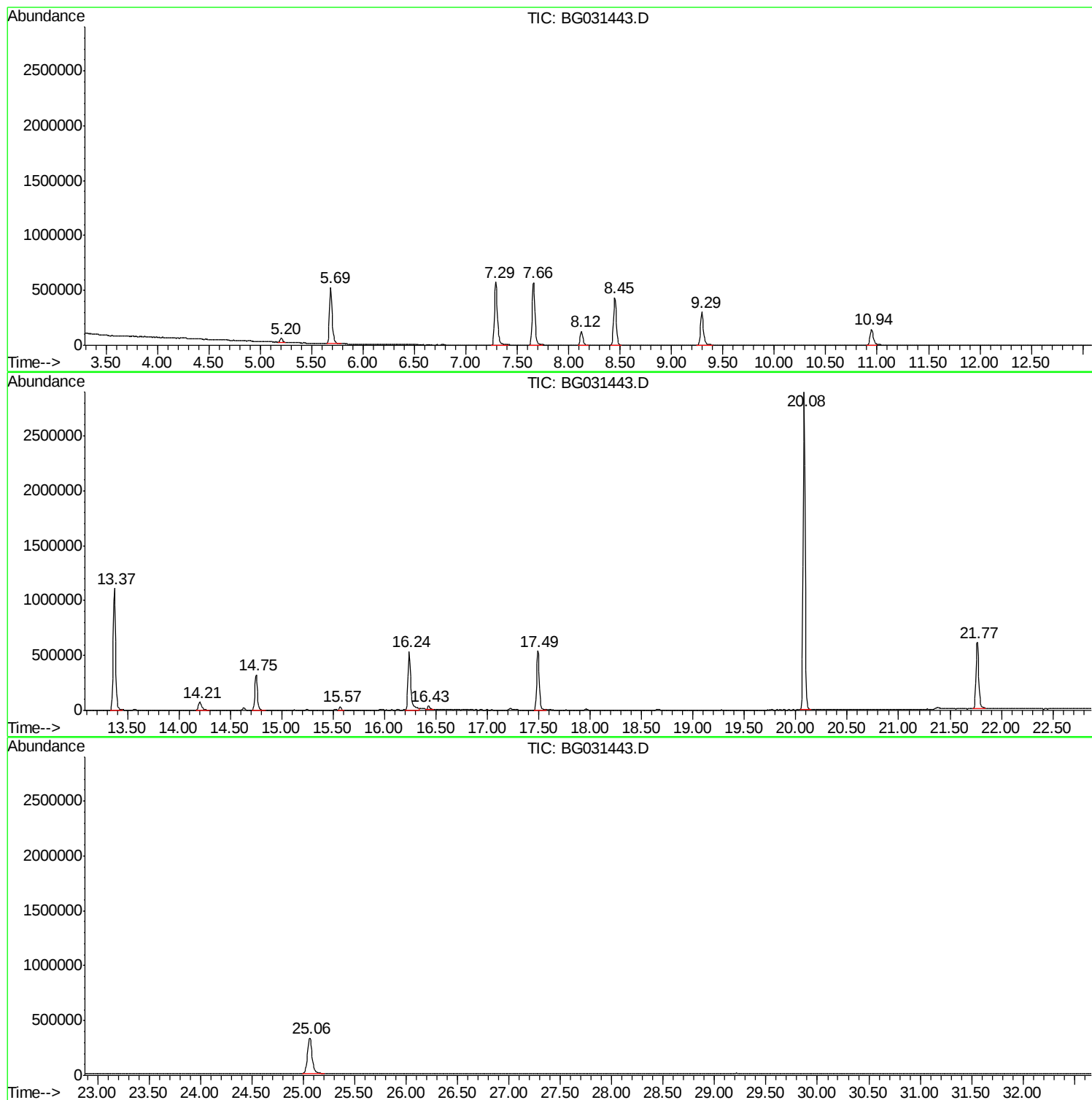
Sum of corrected areas: 15920448

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Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG111017.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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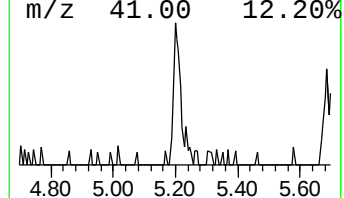
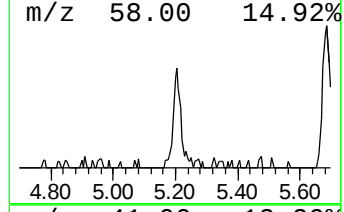
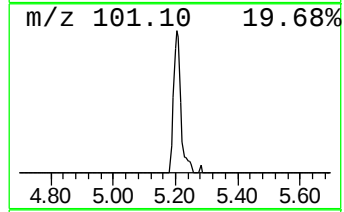
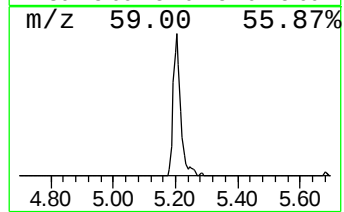
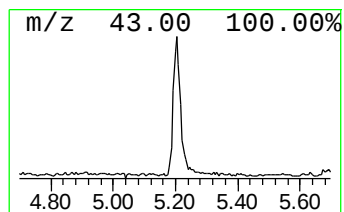
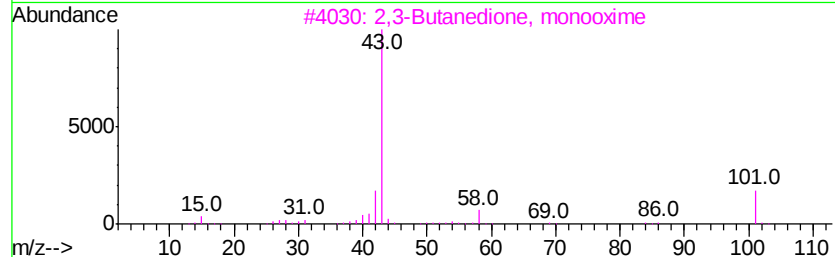
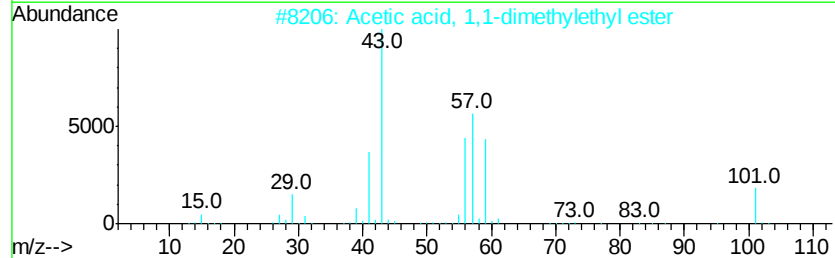
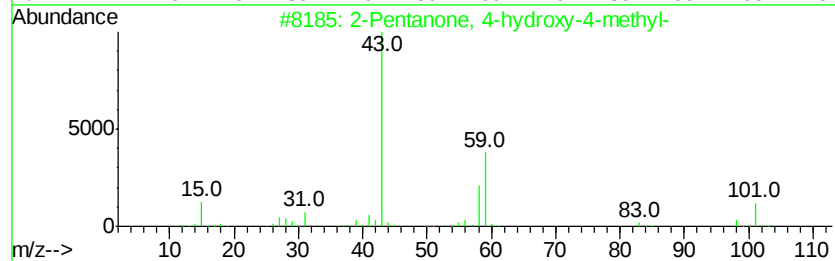
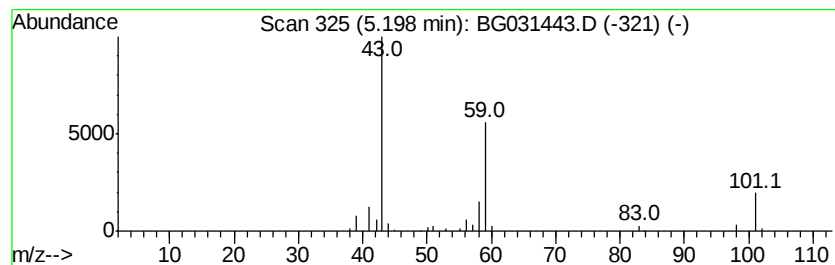
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.20	5.19 ng	58777	1,4-Dichlorobenzene-d4	8.13

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56
2			Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	39
3			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	38
4			Butyl aldoxime, 3-methyl-, syn-	101	C5H11NO	005780-40-5	32
5			2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	28



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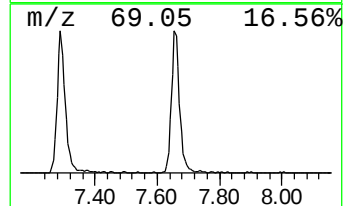
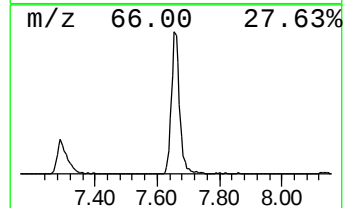
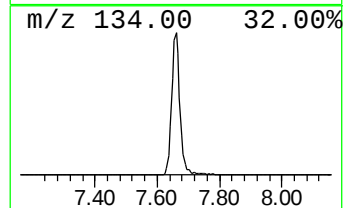
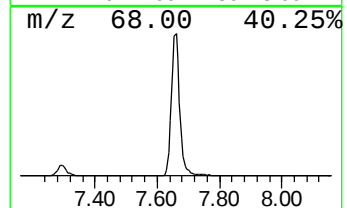
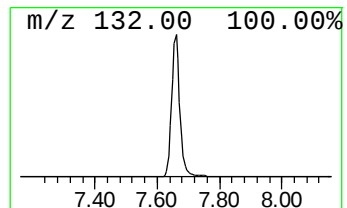
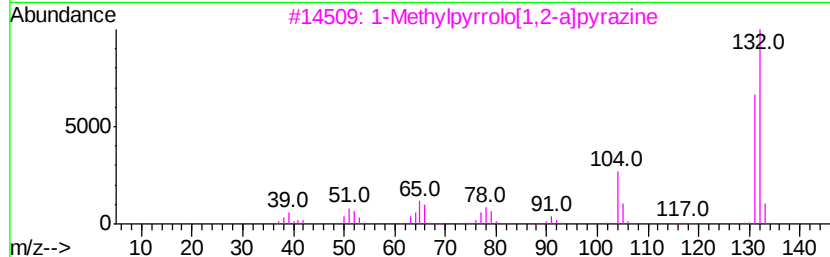
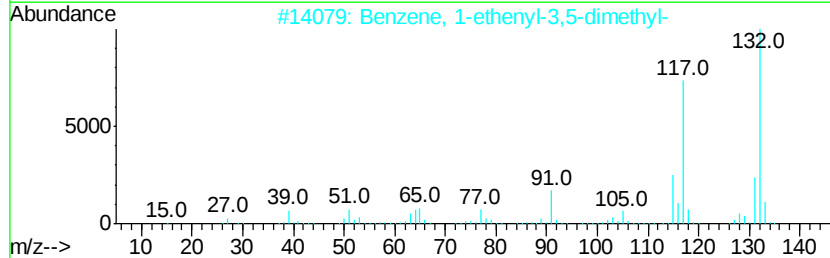
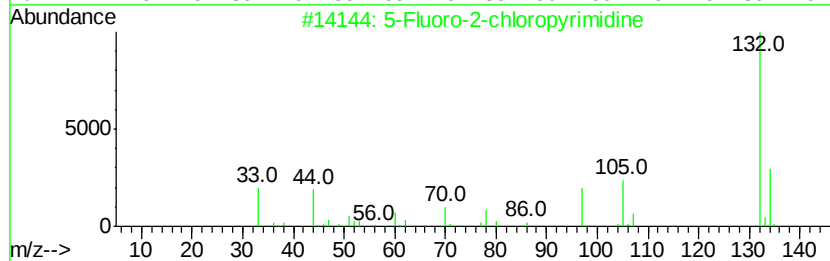
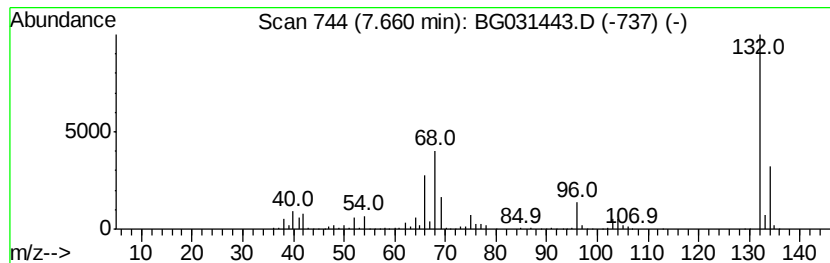
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Peak Number 2 unknown7.66 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.66	93.49 ng	1059330	1,4-Dichlorobenzene-d4	8.13

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	12
2		Benzene, 1-ethenyl-3,5-dimethyl-	132	C10H12	005379-20-4	9
3		1-Methylpyrrolo[1,2-a]pyrazine	132	C8H8N2	064608-59-9	9
4		4-Chloro-2-fluoro-pyrimidine	132	C4H2ClFN2	051422-00-5	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...	5.20	5.2	ng	58777	1	8.13	226614	20.0
unknown7.66	7.66	93.5	ng	1059330	1	8.13	226614	20.0