

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA G\DATA\BG030618\
 Data File : BG033031.D
 Acq On : 6 Mar 2018 21:09
 Operator : SJ/JU
 Sample : PB107125BL
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB107125BL

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-BG022118.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	3.506	32	37	50	rVB5	12123	31808	1.10%	0.193%
2	5.820	423	431	438	rBV	22199	36847	1.27%	0.224%
3	6.326	510	517	533	rBV	861018	1478502	51.02%	8.982%
4	8.006	795	803	817	rBV	835813	1559791	53.83%	9.476%
5	8.423	864	874	887	rBV	823135	1522371	52.53%	9.249%
6	8.910	949	957	965	rVB	179254	318380	10.99%	1.934%
7	9.251	1005	1015	1025	rBV	611045	1136433	39.22%	6.904%
8	10.109	1153	1161	1174	rBV	498349	919644	31.74%	5.587%
9	11.795	1441	1448	1459	rBV	248963	468357	16.16%	2.845%
10	14.156	1842	1850	1859	rBV	1335167	2067008	71.33%	12.557%
11	15.525	2075	2083	2091	rVB2	370806	607695	20.97%	3.692%
12	16.994	2326	2333	2348	rBV	889795	1402262	48.39%	8.519%
13	18.251	2540	2547	2557	rBV	455884	721649	24.90%	4.384%
14	20.771	2970	2976	2987	rBV	2124650	2897839	100.00%	17.605%
15	22.609	3281	3289	3303	rBV	347761	720955	24.88%	4.380%
16	26.410	3924	3936	3961	rBV2	127913	571003	19.70%	3.469%

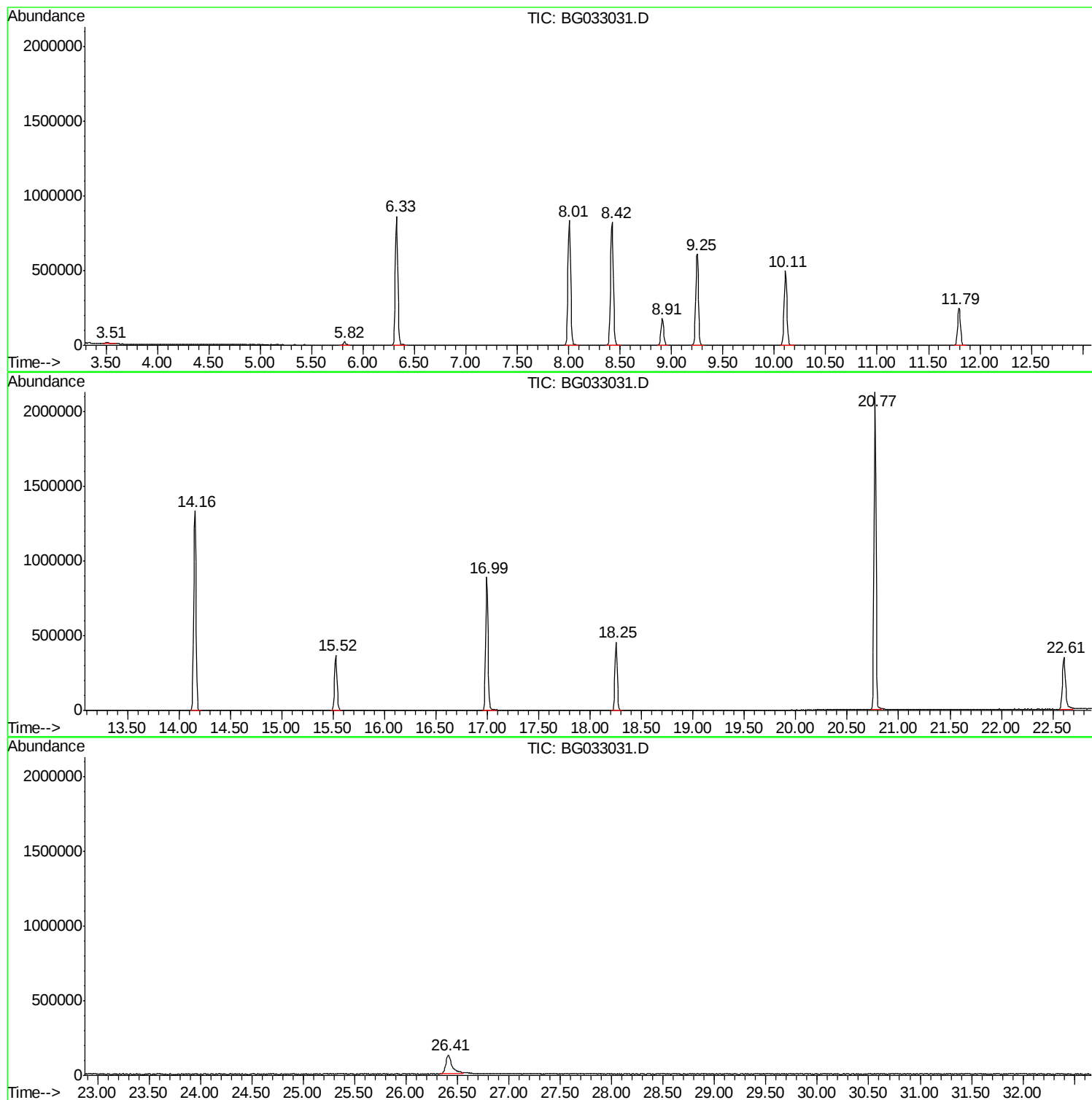
Sum of corrected areas: 16460544

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Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG022118.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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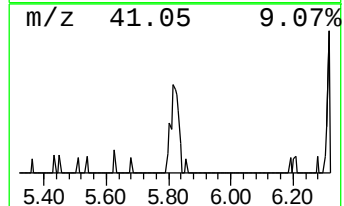
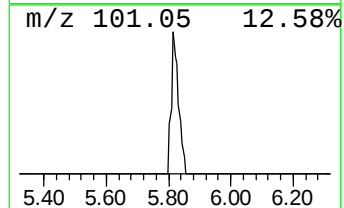
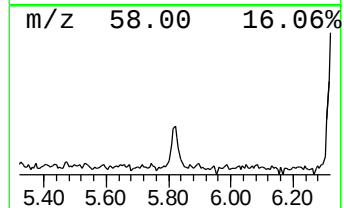
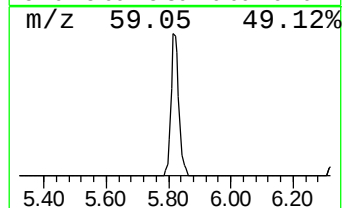
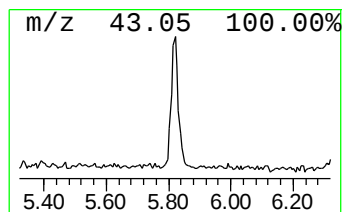
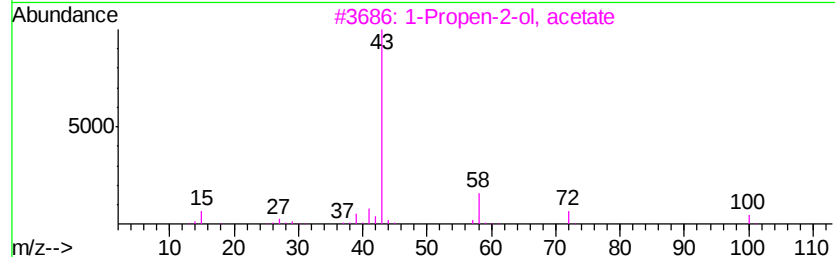
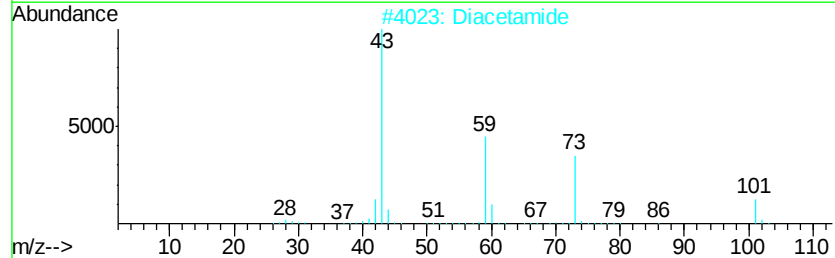
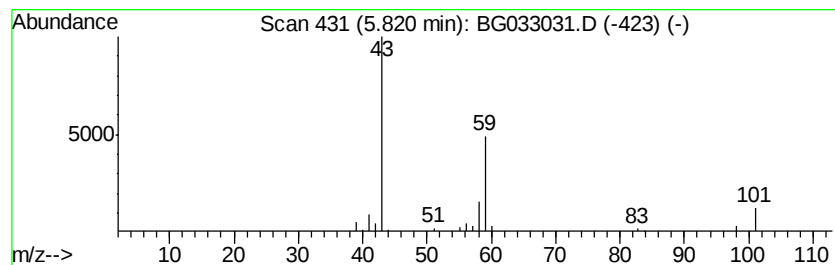
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.	
5.82	2.31 ng	36847	1,4-Dichlorobenzene-d4	8.91	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56
2	Diacetamide	101	C4H7NO2	000625-77-4	9
3	1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	9
4	5-Hexen-2-one	98	C6H10O	000109-49-9	7
5	2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	7



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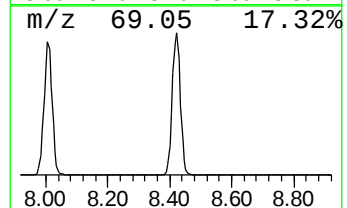
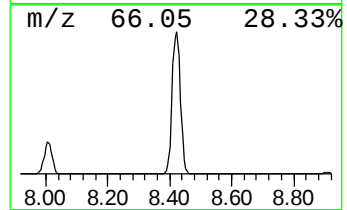
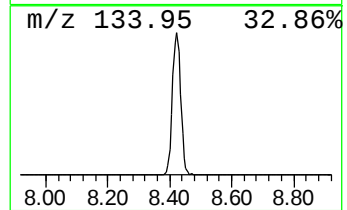
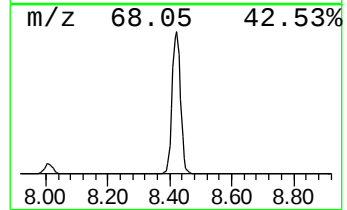
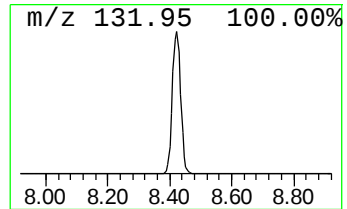
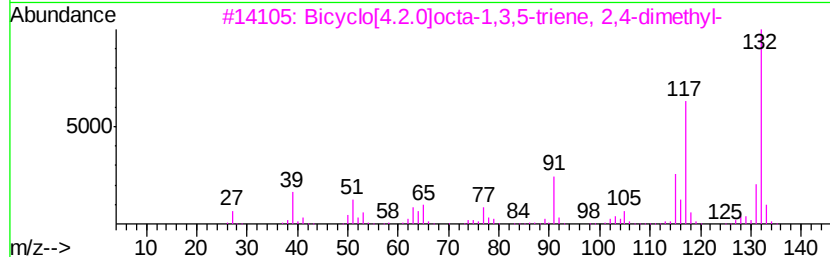
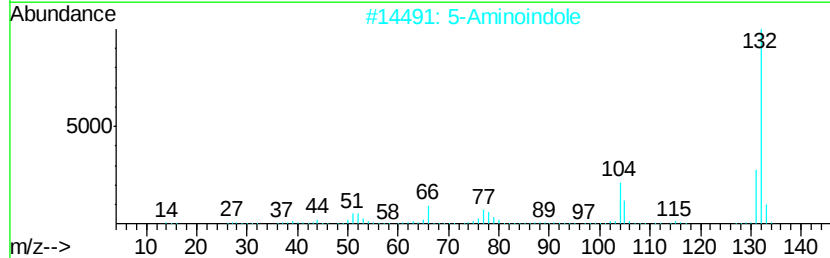
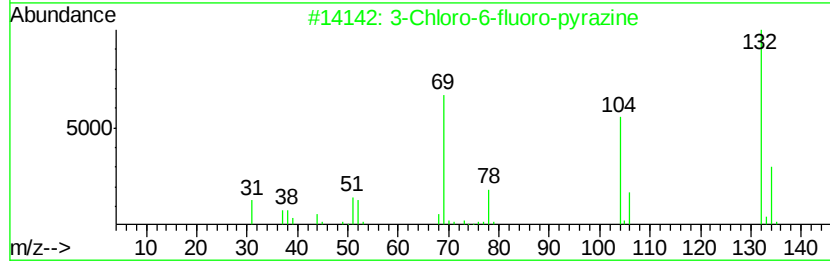
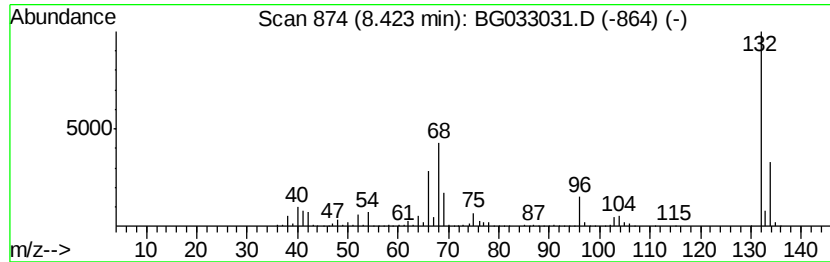
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Peak Number 2 unknown8.42 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.42	95.63 ng	1522370	1,4-Dichlorobenzene-d4	8.91

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	35
2		5-Aminoindole	132	C8H8N2	005192-03-0	12
3		Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9
4		Benzene, 1-ethenyl-3,5-dimethyl-	132	C10H12	005379-20-4	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...	5.82	2.3	ng	36847	1	8.91	318380	20.0
unknown8.42	8.42	95.6	ng	1522370	1	8.91	318380	20.0