

Data Path : Z:\HPCHEM1\BNA G\DATA\BG113017\
 Data File : BG031299.D
 Acq On : 30 Nov 2017 22:15
 Operator : SJ/JU
 Sample : I6634-07
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 112717-TIDO-E-U-B

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	5.206	320	326	335	rBV	64043	104201	1.59%	0.447%
2	5.688	402	408	419	rBV	311214	558909	8.55%	2.396%
3	7.298	675	682	695	rBV	185712	386779	5.91%	1.658%
4	7.662	736	744	759	rBV	623957	1136676	17.38%	4.873%
5	8.132	816	824	832	rBV	222007	389812	5.96%	1.671%
6	8.455	871	879	891	rBV	907408	1659050	25.37%	7.112%
7	9.295	1013	1022	1045	rBV	658612	1295880	19.82%	5.555%
8	10.946	1295	1303	1322	rBV	266749	568969	8.70%	2.439%
9	13.373	1709	1716	1733	rBV	2163673	3732740	57.08%	16.001%
10	14.753	1942	1951	1963	rBV2	519938	924915	14.14%	3.965%
11	16.234	2197	2203	2220	rBV	1115043	1913516	29.26%	8.203%
12	17.491	2411	2417	2433	rBV2	740201	1234810	18.88%	5.293%
13	20.088	2852	2859	2872	rBV	4561430	6539155	100.00%	28.032%
14	21.769	3137	3145	3153	rBV	853480	1478410	22.61%	6.338%
15	25.059	3694	3705	3721	rBV2	466891	1404017	21.47%	6.019%

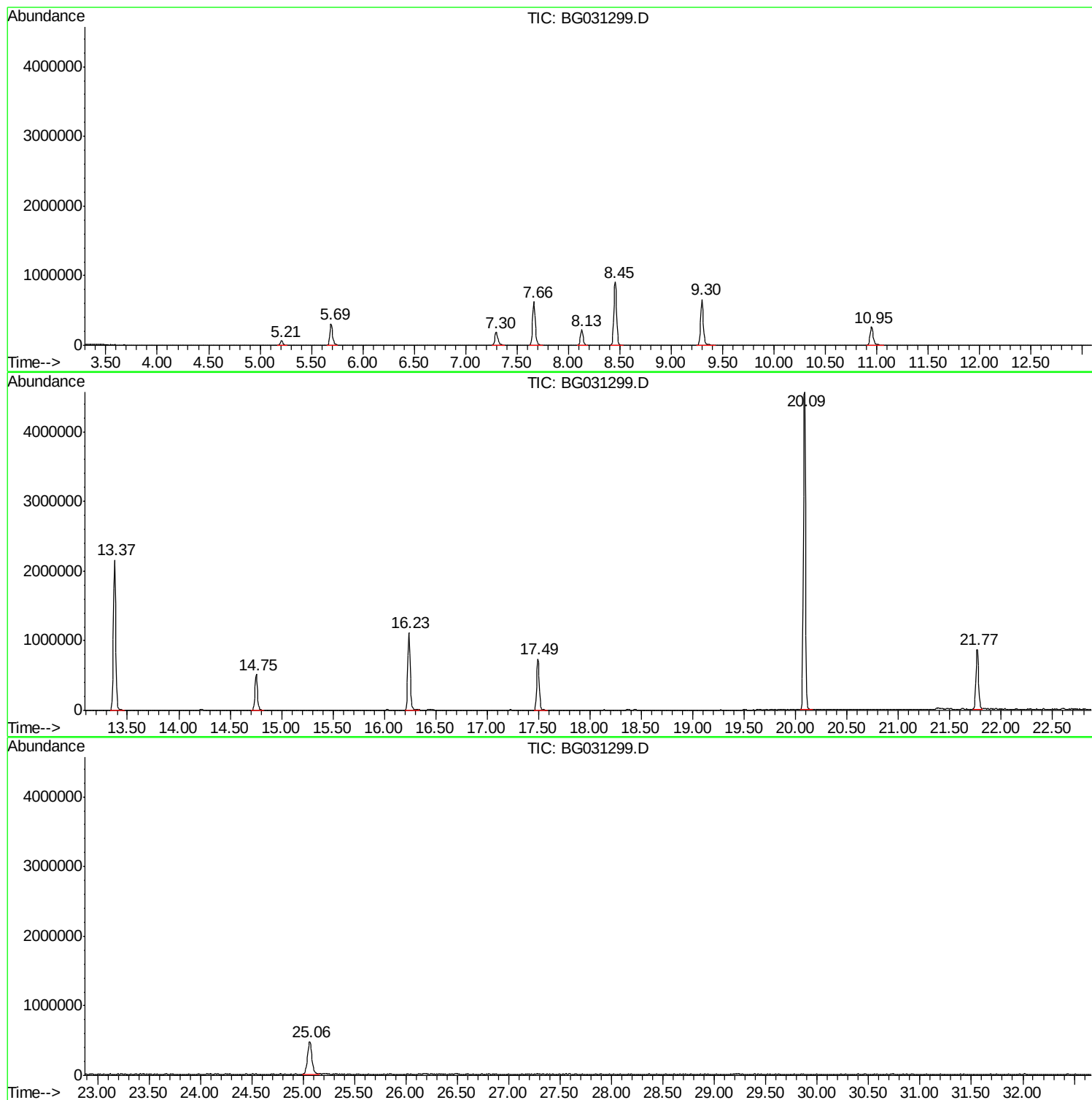
Sum of corrected areas: 23327839

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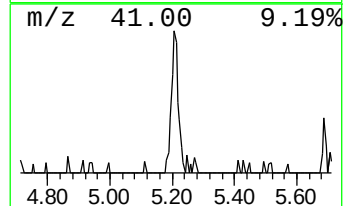
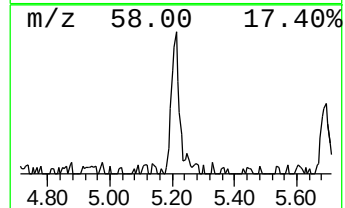
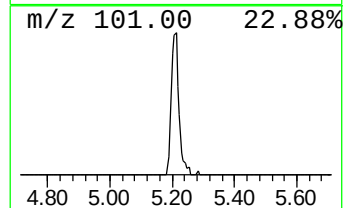
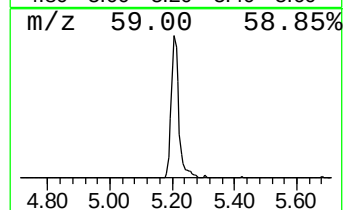
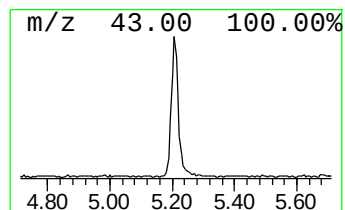
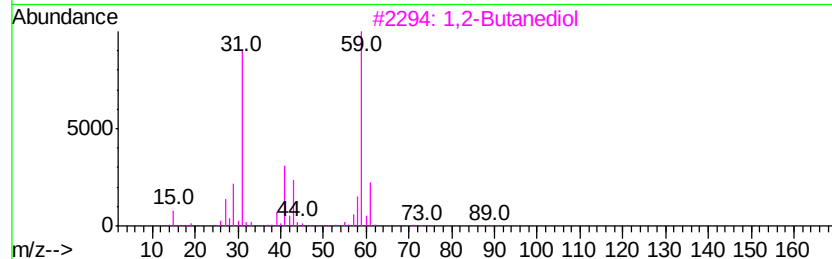
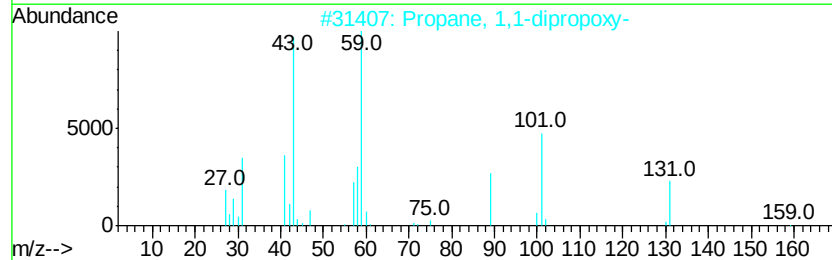
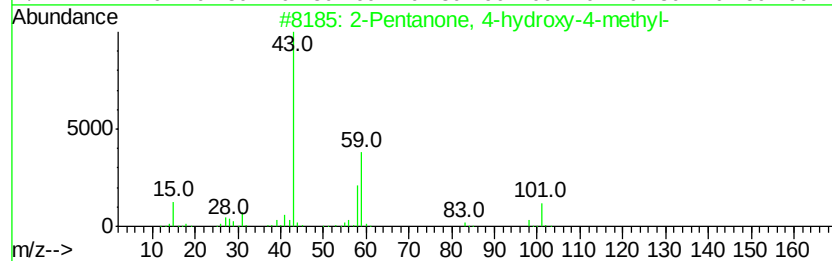
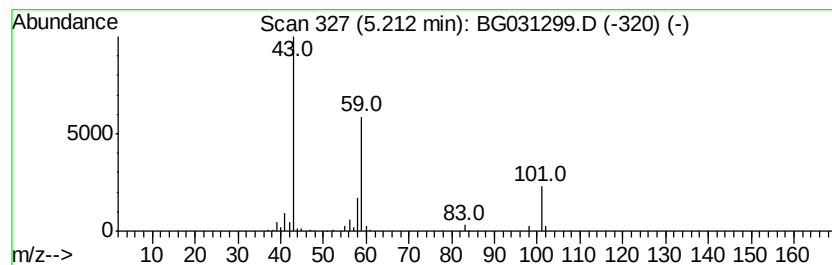
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.21	5.35 ng	104201	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	64
2			Propane, 1,1-dipropoxy-	160	C9H20O2	004744-11-0	38
3			1,2-Butanediol	90	C4H10O2	000584-03-2	23
4			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9
5			Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9



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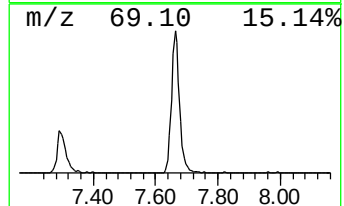
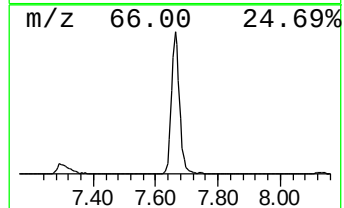
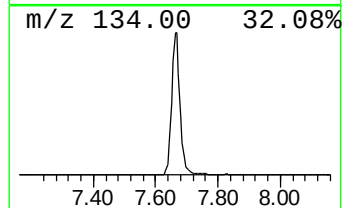
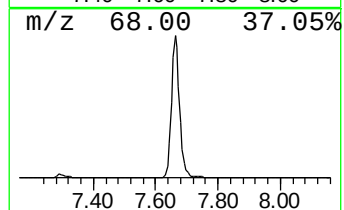
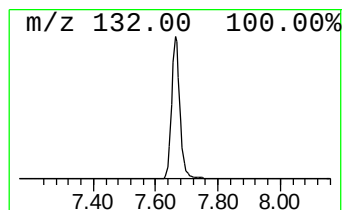
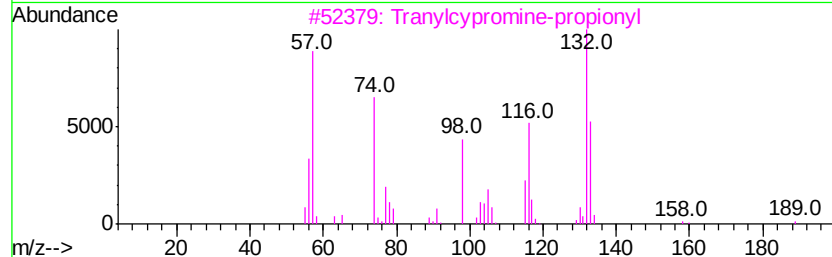
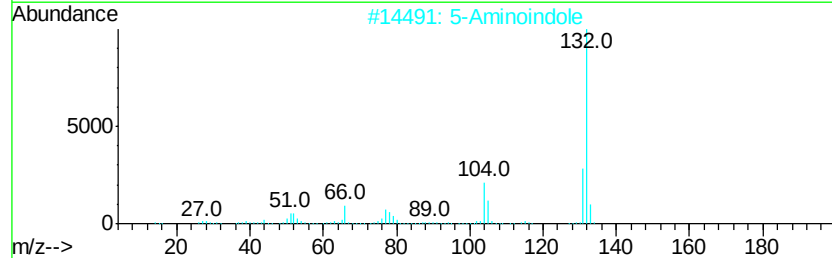
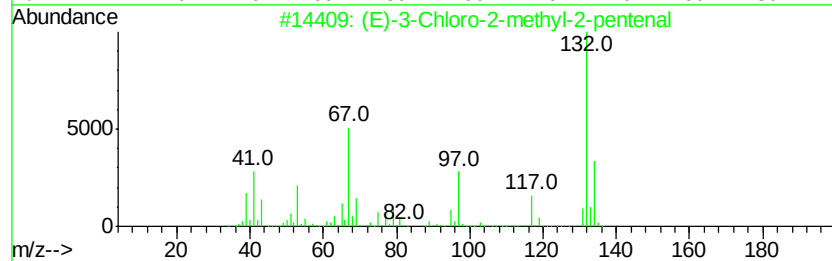
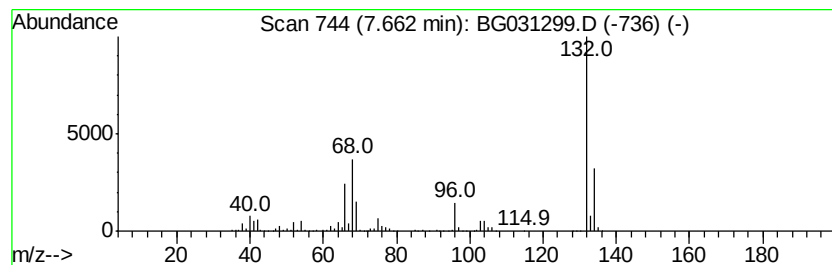
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Peak Number 2 unknown7.66 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.66	58.32 ng	1136680	1,4-Dichlorobenzene-d4	8.13

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	25
2		5-Aminoindole	132	C8H8N2	005192-03-0	12
3		Tranvlcyproline-propionyl	189	C12H15NO	1000123-86-3	12
4		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	9
5		4-Methylpyrrolo[1,2-a]pyrazine	132	C8H8N2	064608-60-2	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.21	5.3	ng	104201	1	8.13	389812	20.0
unknown7.66	7.66	58.3	ng	1136680	1	8.13	389812	20.0