

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG103019\
 Data File : BG043430.D
 Acq On : 31 Oct 2019 11:29
 Operator : JU
 Sample : PB124413BL
 Misc :
 ALS Vial : 29 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB124413BL

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:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG102119.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	5.124	341	346	357	rBV	125748	192079	5.27%	1.218%
2	5.612	420	429	444	rBV	627144	1114254	30.54%	7.064%
3	7.210	691	701	712	rBV	640229	1212158	33.23%	7.684%
4	7.562	753	761	781	rBV	701349	1235985	33.88%	7.835%
5	8.021	832	839	853	rVB2	126182	225493	6.18%	1.429%
6	8.344	887	894	904	rBV	520166	944246	25.88%	5.986%
7	9.184	1028	1037	1052	rBV	340993	707837	19.40%	4.487%
8	10.829	1309	1317	1325	rBV	128082	259457	7.11%	1.645%
9	13.273	1726	1733	1747	rBV	1147343	1907452	52.29%	12.092%
10	14.648	1960	1967	1979	rBV2	251778	450592	12.35%	2.856%
11	16.140	2213	2221	2239	rBV	896225	1714989	47.01%	10.872%
12	17.398	2428	2435	2449	rBV	309924	583423	15.99%	3.698%
13	20.006	2872	2879	2896	rBV	2474553	3648174	100.00%	23.127%
14	21.663	3153	3161	3178	rBV	375124	765677	20.99%	4.854%
15	24.860	3693	3705	3722	rBV3	266223	812966	22.28%	5.154%

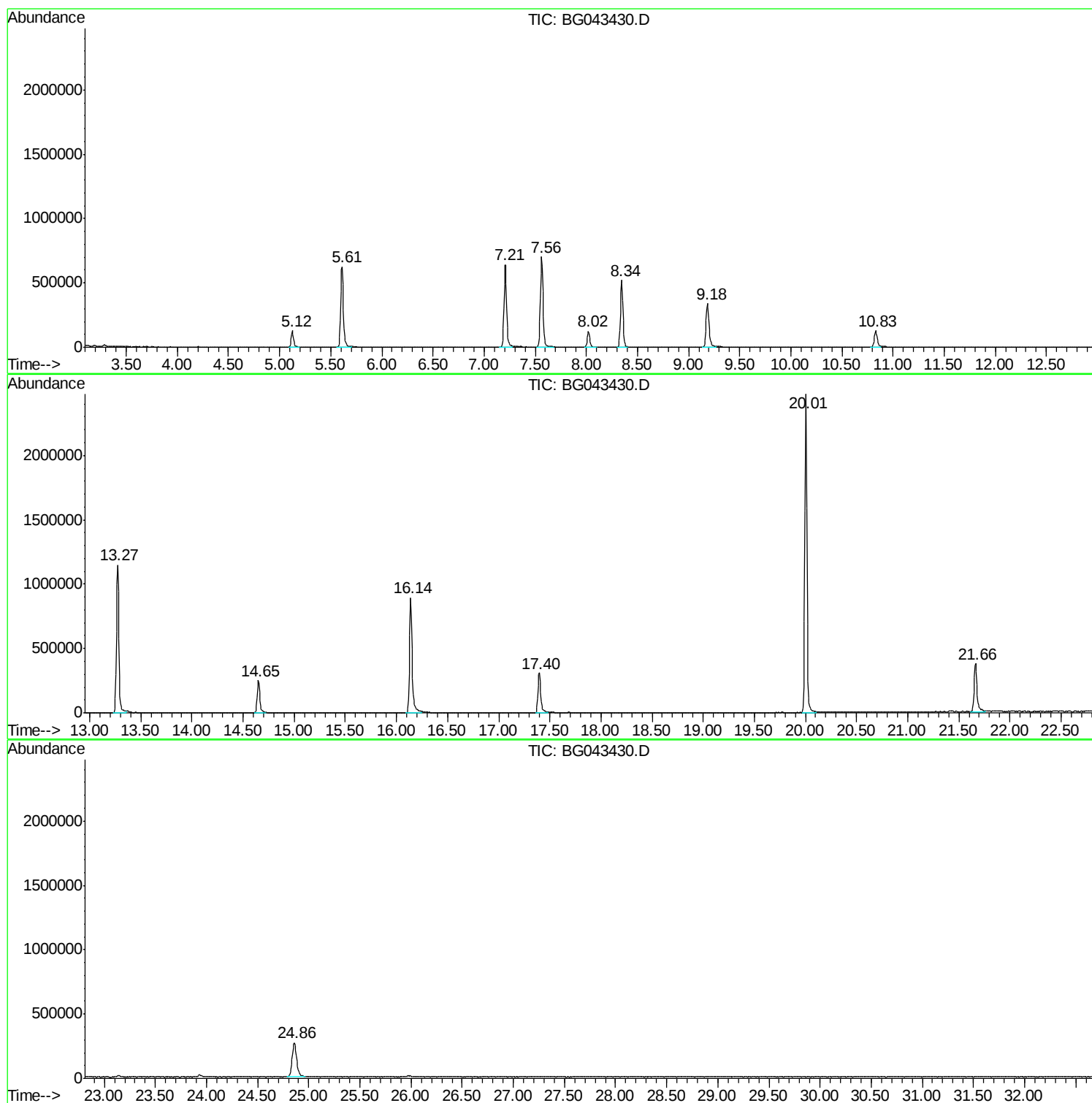
Sum of corrected areas: 15774782

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Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG102119.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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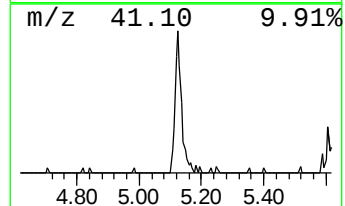
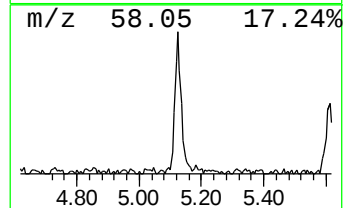
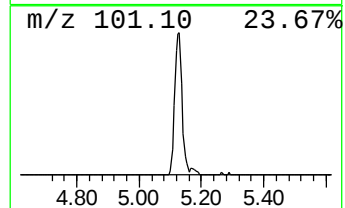
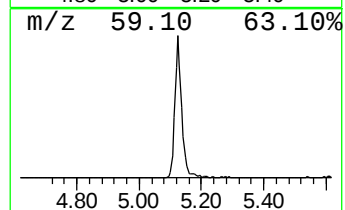
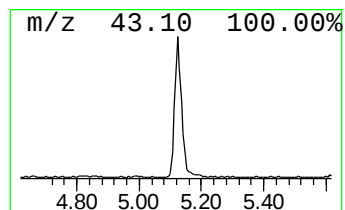
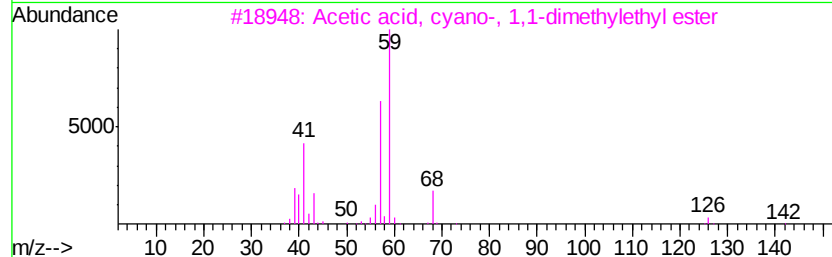
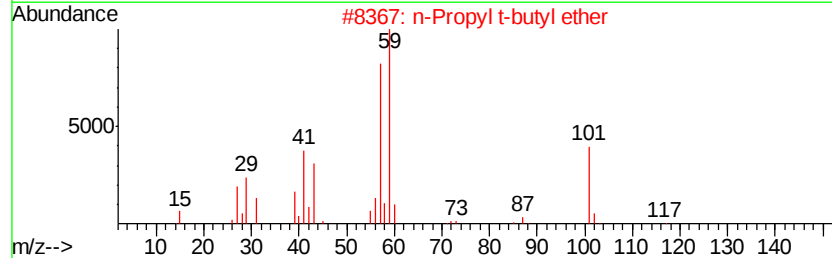
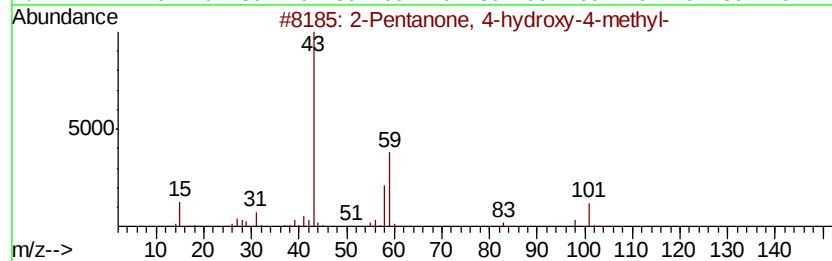
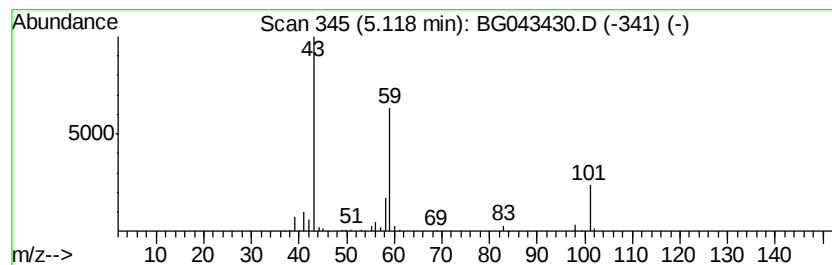
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.12	17.04 ng	192079	1,4-Dichlorobenzene-d4	8.02

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56
2			n-Propyl t-butyl ether	116	C7H16O	1000334-81-9	33
3			Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	32
4			Butyl aldoxime, 3-methyl-, syn-	101	C5H11NO	005780-40-5	17
5			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	16



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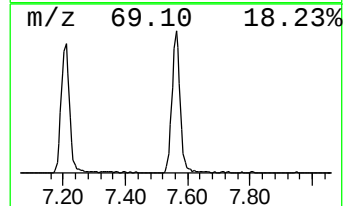
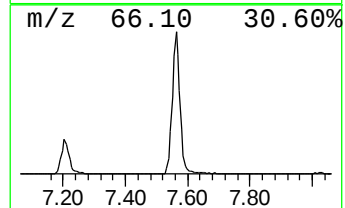
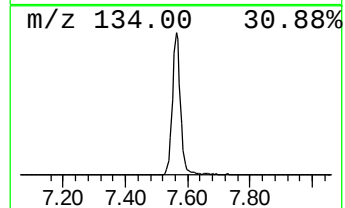
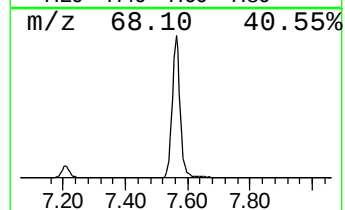
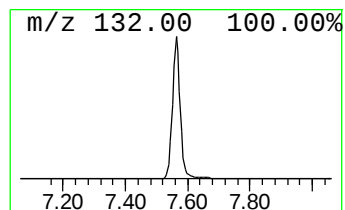
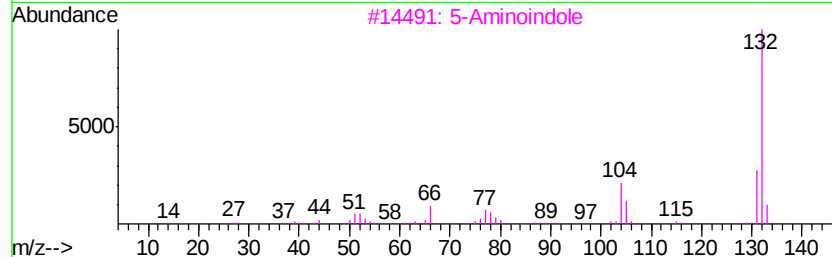
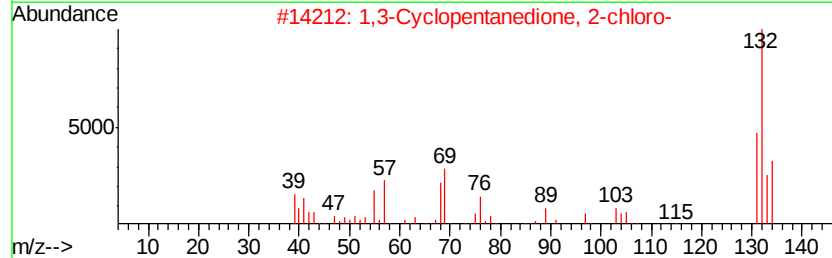
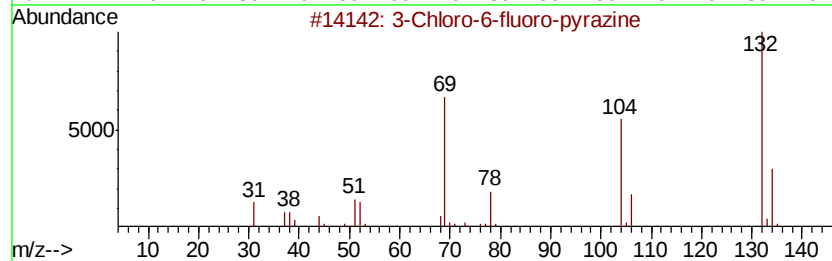
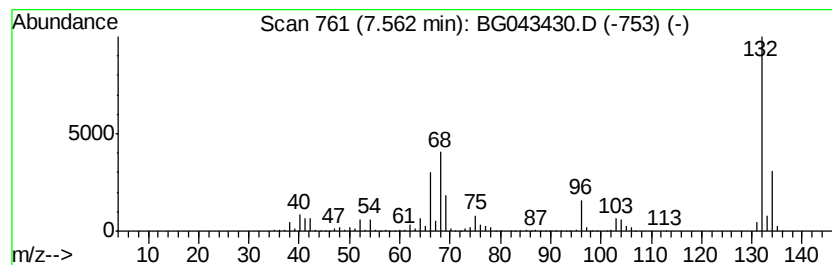
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Peak Number 2 unknown7.56 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.56	109.63 ng	1235990	1,4-Dichlorobenzene-d4	8.02

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		1,3-Cyclopentanedione, 2-chloro-	132	C5H5ClO2	014203-19-1	17
3		5-Aminoindole	132	C8H8N2	005192-03-0	10
4		(5-Methyl-2-pyridyl)acetonitrile	132	C8H8N2	1000241-93-9	10
5		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.12	17.0	ng	192079	1	8.02	225493	20.0
unknown7.56	7.56	109.6	ng	1235990	1	8.02	225493	20.0