

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG010220\
 Data File : BG043871.D
 Acq On : 2 Jan 2020 12:03
 Operator : JU
 Sample : PB125838BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB125838BL

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-BG123019.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	5.068	331	337	350	rVB	635180	978028	8.80%	1.677%
2	5.539	409	417	430	rBV	2791384	4473444	40.23%	7.673%
3	7.131	678	688	700	rBV	2837350	5028791	45.23%	8.625%
4	7.495	742	750	761	rBV	2790236	4973316	44.73%	8.530%
5	7.965	822	830	838	rVB	520654	872822	7.85%	1.497%
6	8.288	877	885	893	rBV	2176939	3734051	33.58%	6.405%
7	9.117	1018	1026	1039	rBV	1691240	3144450	28.28%	5.393%
8	10.768	1299	1307	1319	rVB	711869	1293606	11.63%	2.219%
9	13.224	1716	1725	1734	rBV	4401569	7245871	65.17%	12.428%
10	14.593	1950	1958	1967	rBV2	1183752	1909004	17.17%	3.274%
11	16.079	2204	2211	2235	rBV	3839331	6037061	54.30%	10.355%
12	17.336	2418	2425	2436	rBV	1635744	2376939	21.38%	4.077%
13	19.957	2864	2871	2880	rBV	7712481	11118408	100.00%	19.070%
14	21.590	3142	3149	3159	rVB	1554982	2560233	23.03%	4.391%
15	24.716	3671	3681	3692	rBV2	936053	2556752	23.00%	4.385%

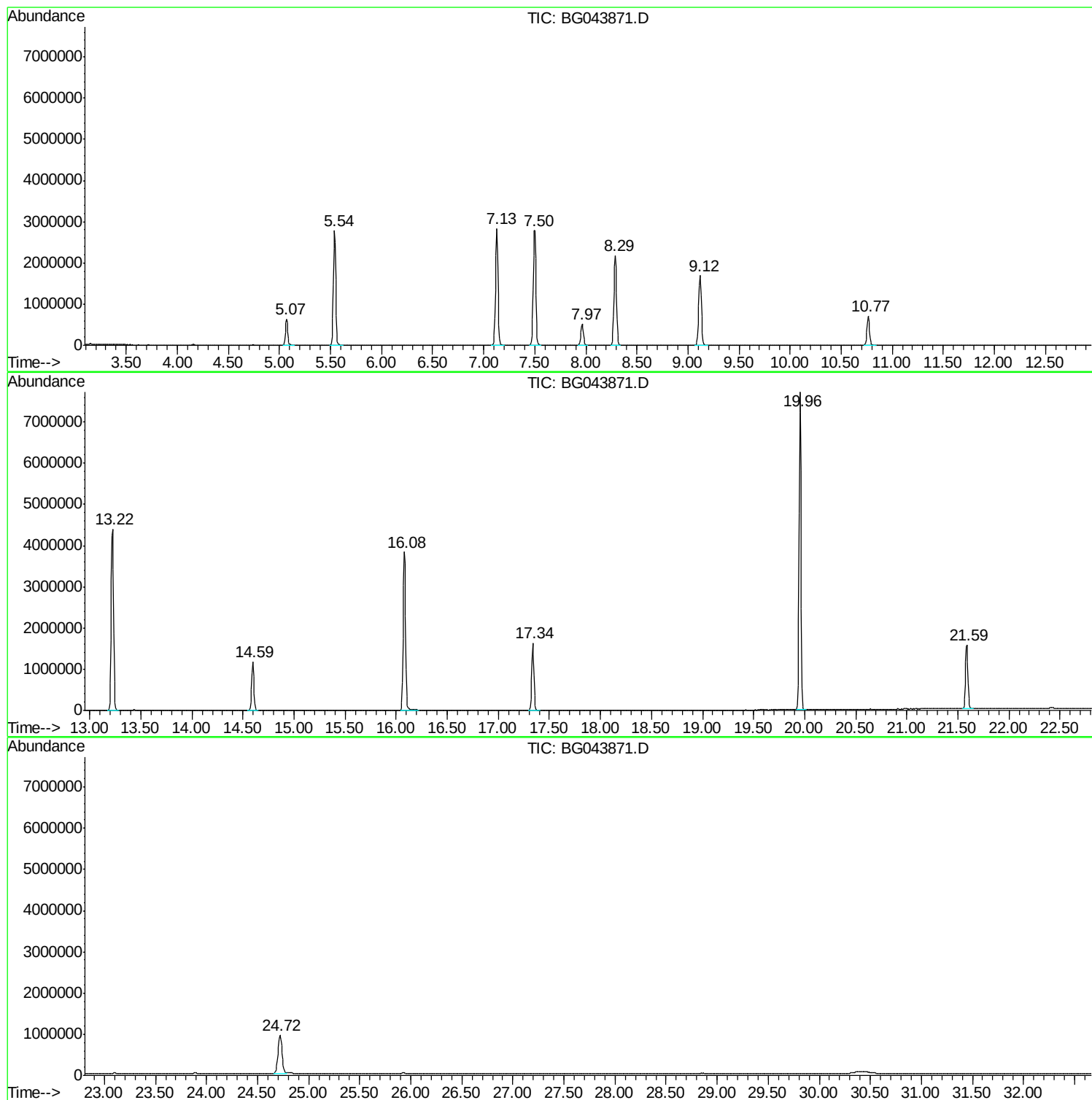
Sum of corrected areas: 58302776

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Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG123019.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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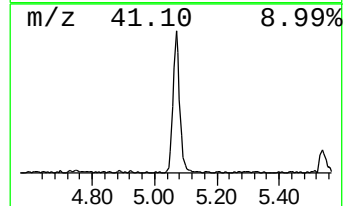
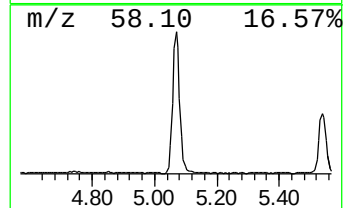
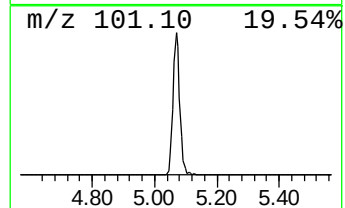
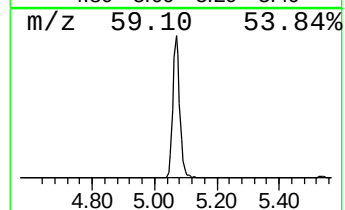
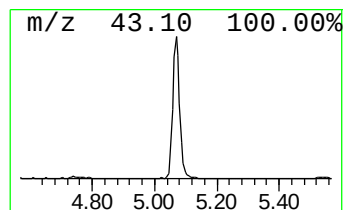
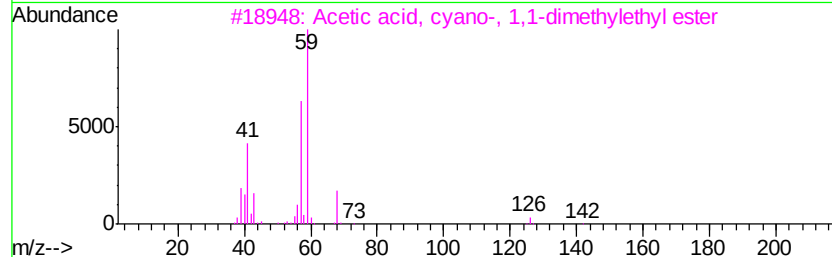
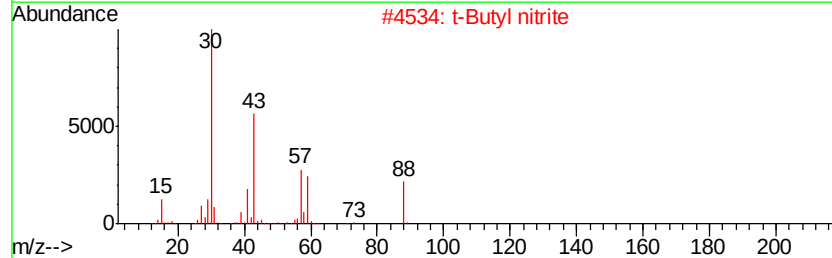
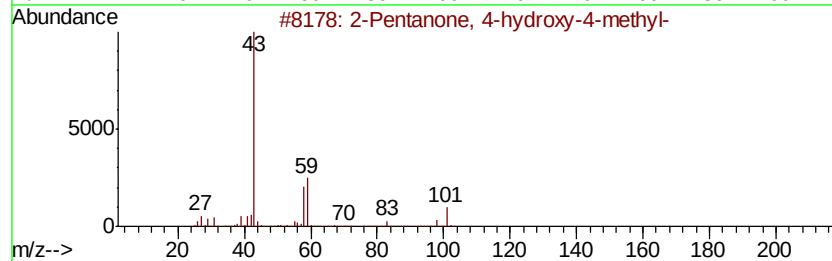
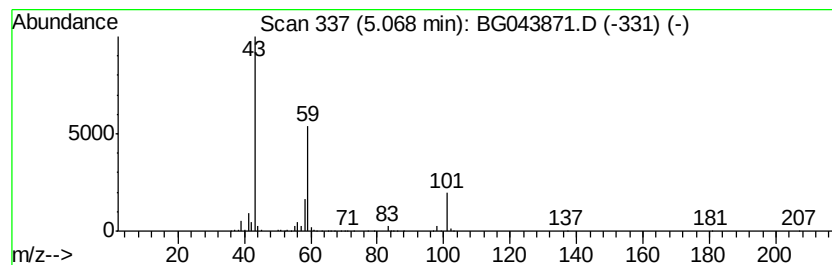
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.07	22.41 ng	978028	1,4-Dichlorobenzene-d4	7.97

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56
2		t-Butyl nitrite	103	C4H9NO2	000540-80-7	45
3		Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	17
4		2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	12
5		Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9



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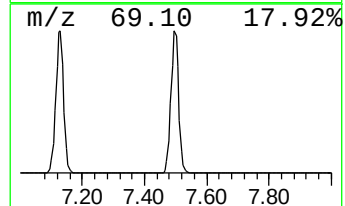
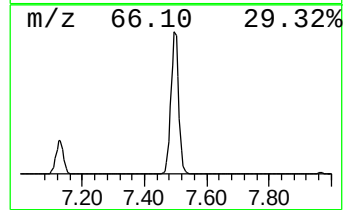
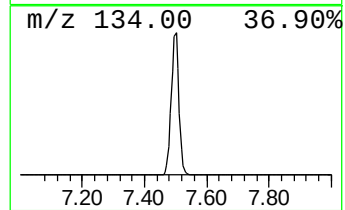
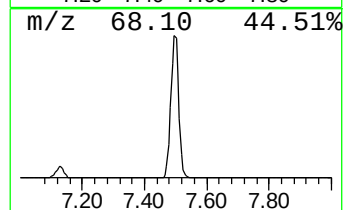
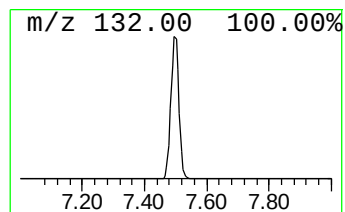
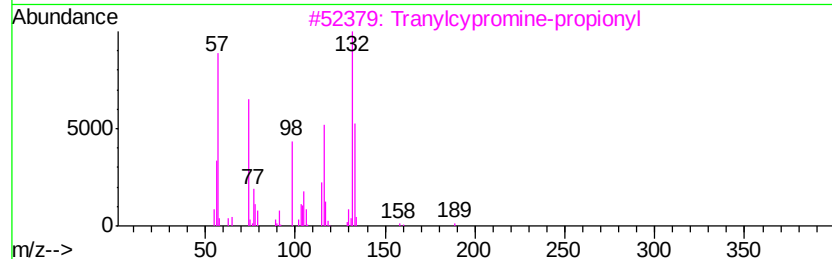
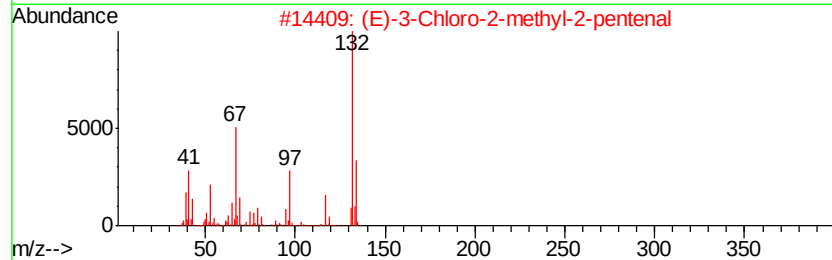
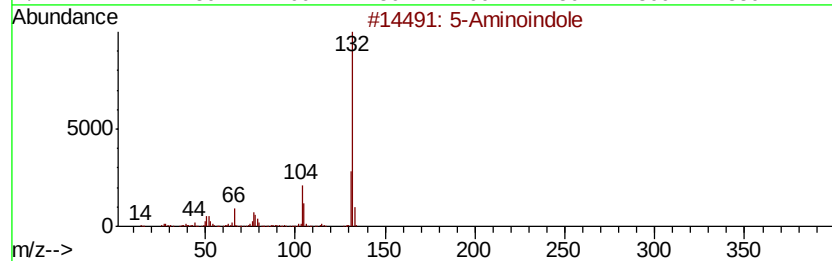
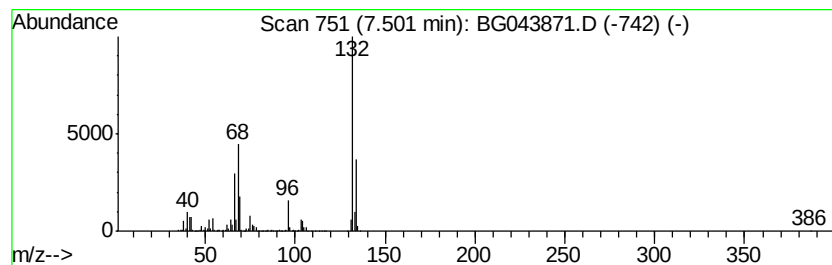
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Peak Number 2 unknown7.50 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.50	113.96 ng	4973320	1,4-Dichlorobenzene-d4	7.97

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	5-Aminoindole	132	C8H8N2	005192-03-0	22
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	12
3		Tranvlcypropine-propionyl	189	C12H15NO	1000123-86-3	12
4		1-Isoquinolinemethanol, .alpha.-...	191	C12H17NO	114608-92-3	12
5		Amphetaminil	250	C17H18N2	017590-01-1	12



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.07	22.4	ng	978028	1	7.97	872822	20.0
unknown7.50	7.50	114.0	ng	4973320	1	7.97	872822	20.0