

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG092518\
 Data File : BG036962.D
 Acq On : 25 Sep 2018 23:29
 Operator : JU/SJ
 Sample : PB113203BL
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB113203BL

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-BG092018.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	5.148	311	316	324	rVB2	31104	53257	1.44%	0.305%
2	5.618	389	396	417	rBV	739390	1354661	36.72%	7.759%
3	7.204	658	666	690	rBV	743647	1484929	40.26%	8.506%
4	7.574	722	729	752	rBV	721395	1382436	37.48%	7.918%
5	8.044	802	809	819	rBV	133439	258312	7.00%	1.480%
6	8.361	856	863	886	rBV	554472	1066043	28.90%	6.106%
7	9.202	999	1006	1031	rBV	427398	921374	24.98%	5.278%
8	10.847	1279	1286	1303	rBV	178912	361063	9.79%	2.068%
9	13.291	1695	1702	1722	rBV	1249643	2107386	57.13%	12.071%
10	14.666	1930	1936	1955	rBV2	300927	543948	14.75%	3.116%
11	16.152	2183	2189	2212	rBV	993539	1762101	47.77%	10.093%
12	17.410	2397	2403	2417	rBV2	423933	737743	20.00%	4.226%
13	20.018	2841	2847	2862	rBV	2725511	3688768	100.00%	21.129%
14	21.675	3123	3129	3147	rVB2	510271	898640	24.36%	5.147%
15	24.883	3665	3675	3686	rBV2	279687	837664	22.71%	4.798%

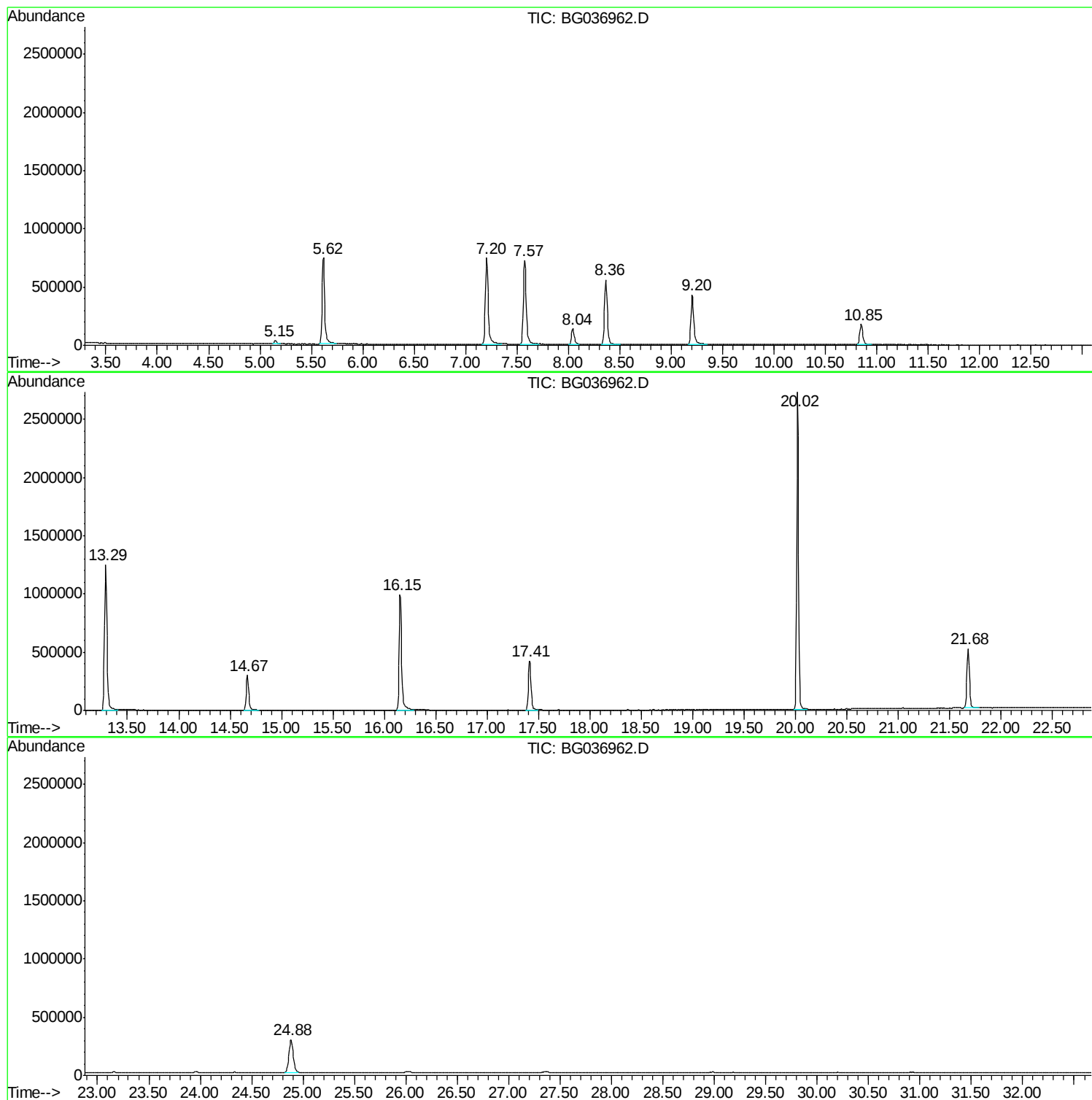
Sum of corrected areas: 17458325

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Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG092018.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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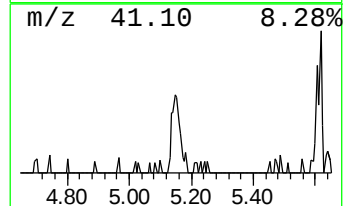
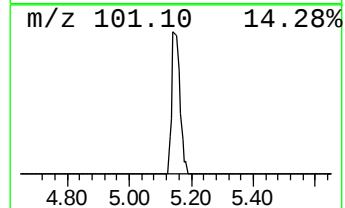
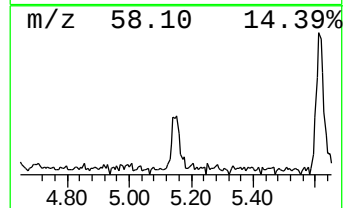
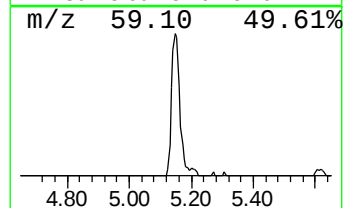
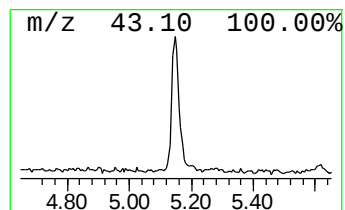
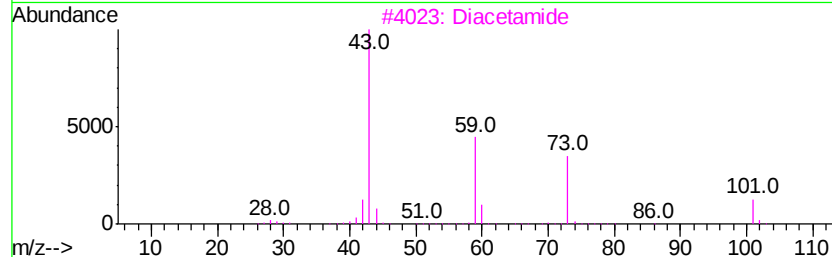
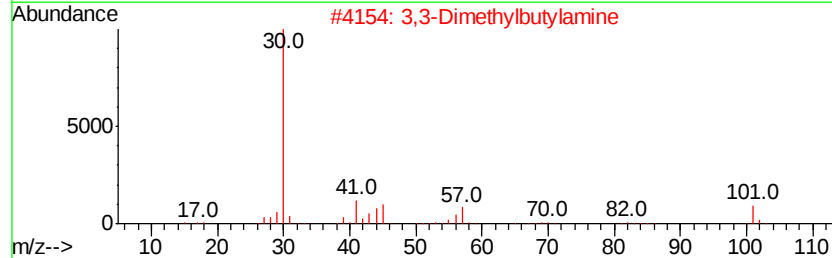
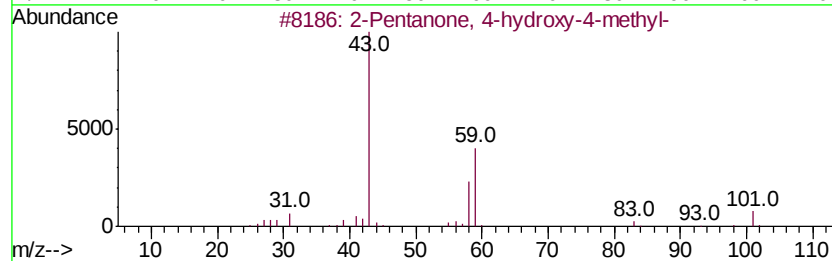
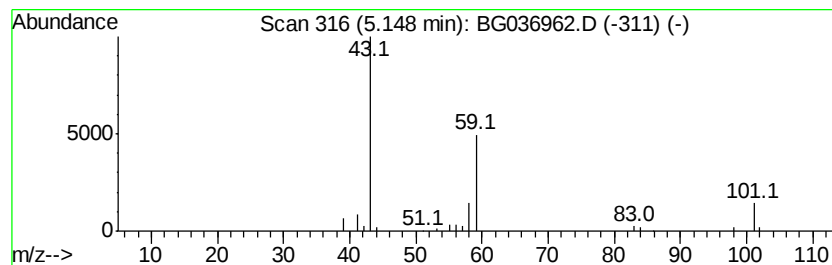
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.15	4.12 ng	53257	1,4-Dichlorobenzene-d4	8.04

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	39
2		3,3-Dimethylbutylamine	101	C6H15N	015673-00-4	9
3		Diacetamide	101	C4H7NO2	000625-77-4	9
4		4-Penten-2-one, 4-methyl-	98	C6H10O	003744-02-3	9
5		Formamide, N,N-diethyl-	101	C5H11NO	000617-84-5	7



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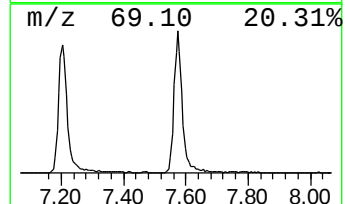
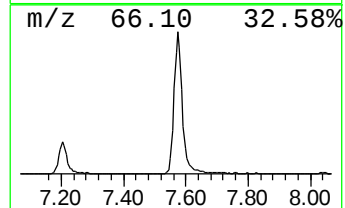
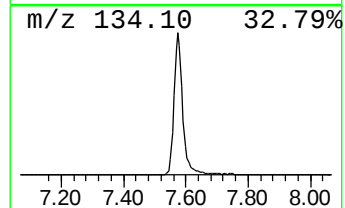
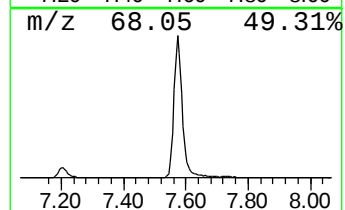
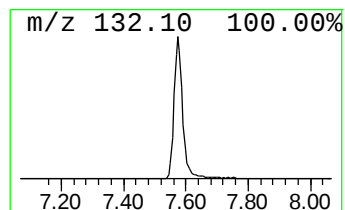
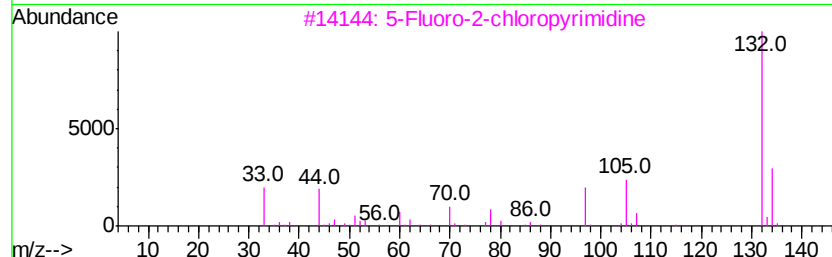
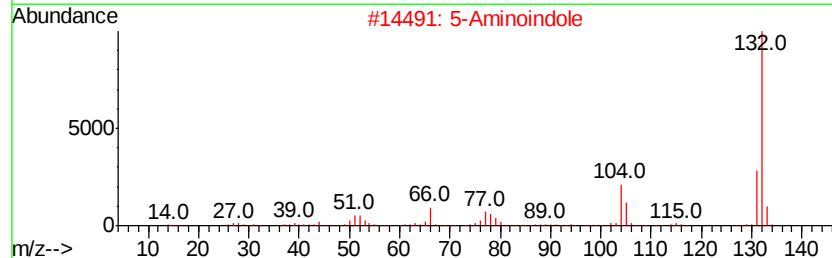
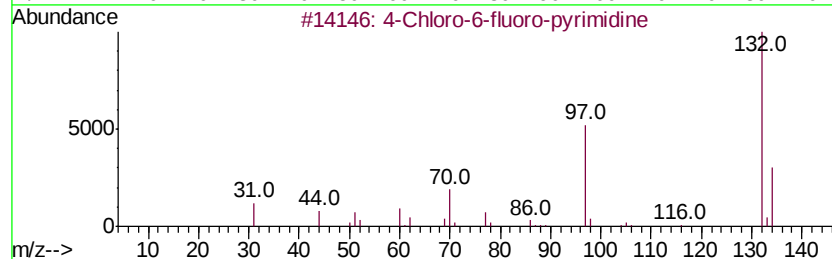
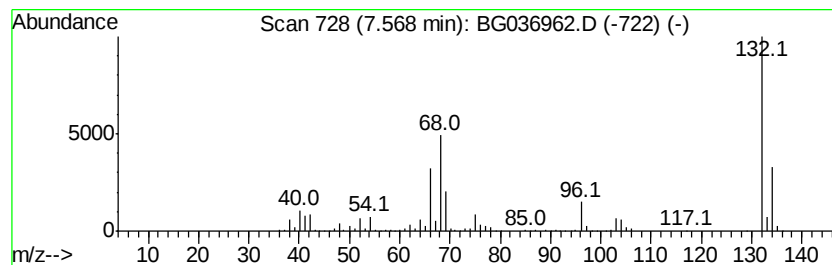
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Peak Number 2 unknown7.57 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.57	107.04 ng	1382440	1,4-Dichlorobenzene-d4	8.04

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	4-Chloro-6-fluoro-pyrimidine	132	C4H2ClFN2	051422-01-6	27
2		5-Aminoindole	132	C8H8N2	005192-03-0	22
3		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	12
4		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
5		(5-Methyl-2-pyridyl)acetonitrile	132	C8H8N2	1000241-93-9	10



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.15	4.1	ng	53257	1	8.04	258312	20.0
unknown7.57	7.57	107.0	ng	1382440	1	8.04	258312	20.0