

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG102918\
 Data File : BG037642.D
 Acq On : 30 Oct 2018 3:37
 Operator : JU/SJ
 Sample : PB114209BL
 Misc :
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB114209BL

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-BG101818.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	5.179	317	322	330	rVB	32467	52336	1.67%	0.313%
2	5.643	394	401	410	rBV	735646	1208179	38.54%	7.217%
3	7.247	665	674	683	rBV	778286	1383776	44.14%	8.266%
4	7.623	730	738	751	rBV	705341	1288114	41.09%	7.695%
5	8.099	812	819	826	rVB	168262	301458	9.62%	1.801%
6	8.422	866	874	884	rVB	528147	947300	30.22%	5.659%
7	9.274	1009	1019	1028	rBV	432692	822069	26.22%	4.911%
8	10.919	1291	1299	1308	rBV	245023	460561	14.69%	2.751%
9	13.358	1706	1714	1723	rBV	1313257	2114254	67.44%	12.630%
10	14.733	1940	1948	1958	rBV2	430128	725259	23.13%	4.332%
11	16.219	2194	2201	2215	rBV	938036	1538101	49.06%	9.188%
12	17.482	2409	2416	2423	rBV	605214	924481	29.49%	5.523%
13	20.079	2852	2858	2869	rBV	2358891	3135163	100.00%	18.728%
14	21.765	3138	3145	3153	rBV2	547527	970062	30.94%	5.795%
15	25.103	3700	3713	3731	rBV3	243673	869089	27.72%	5.192%

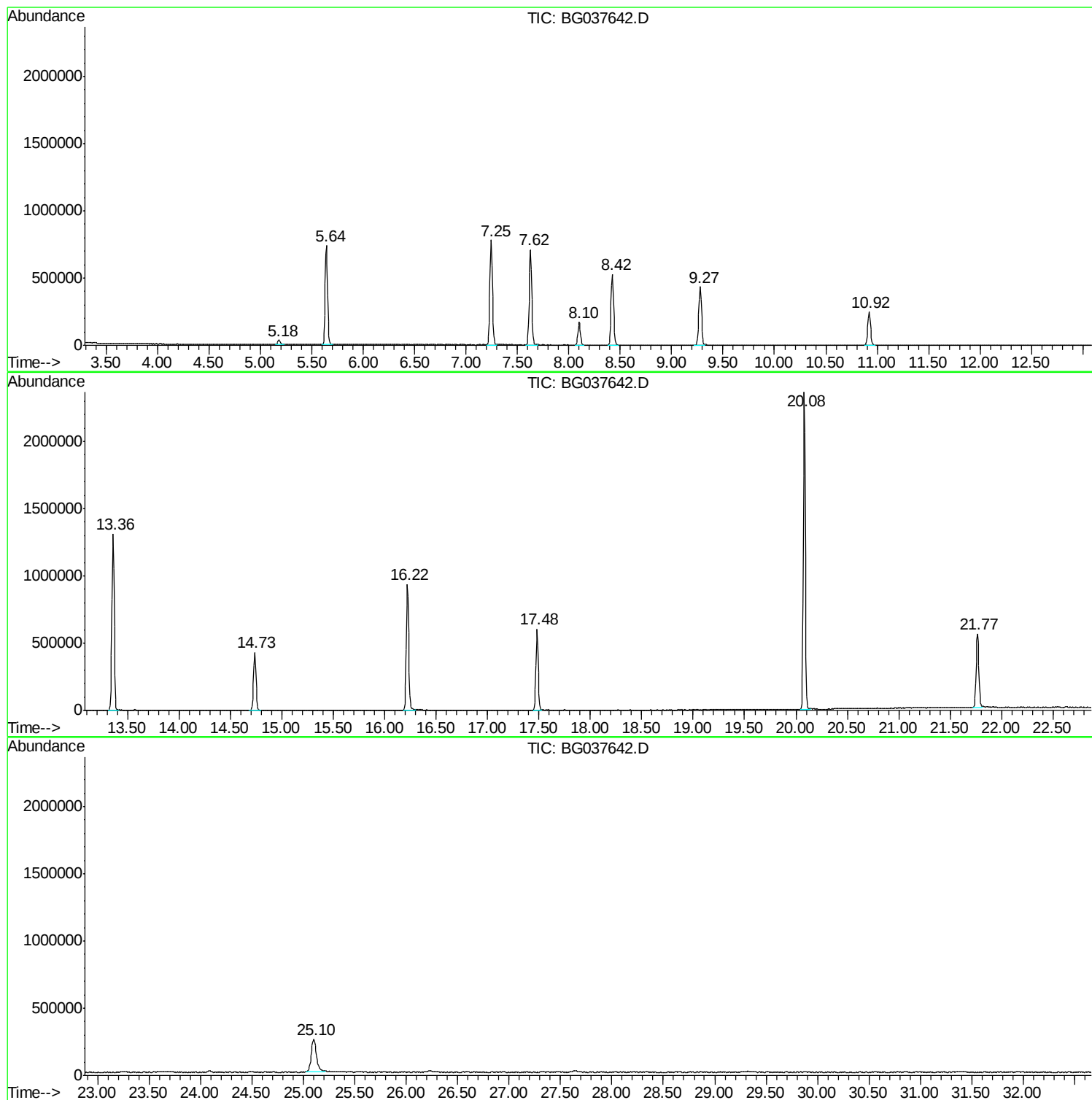
Sum of corrected areas: 16740202

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TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P



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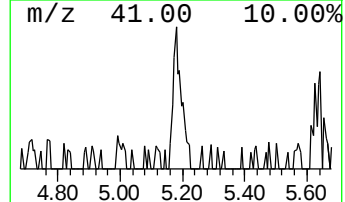
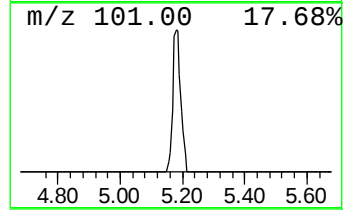
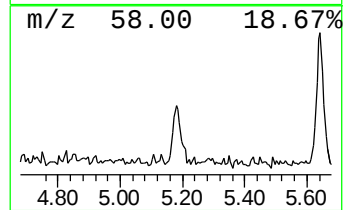
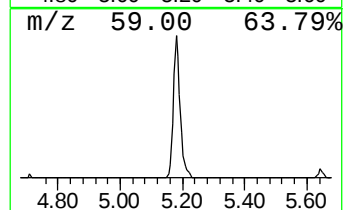
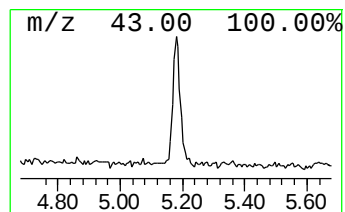
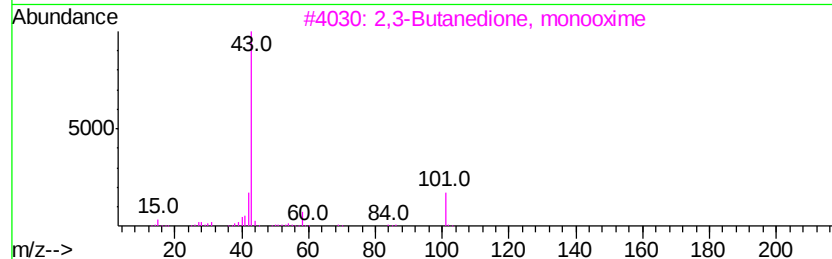
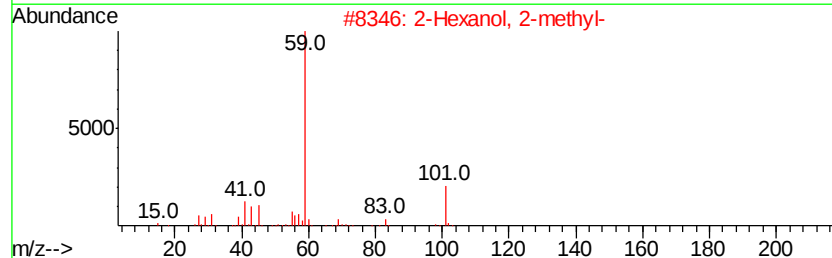
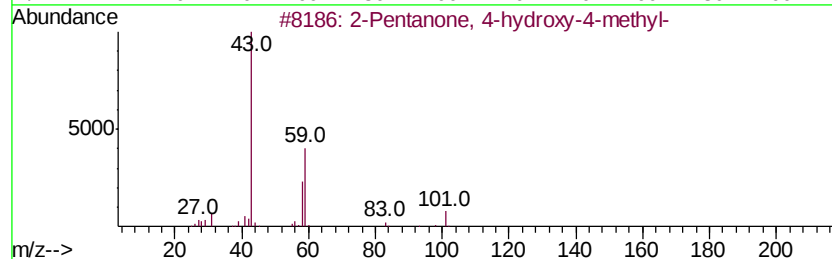
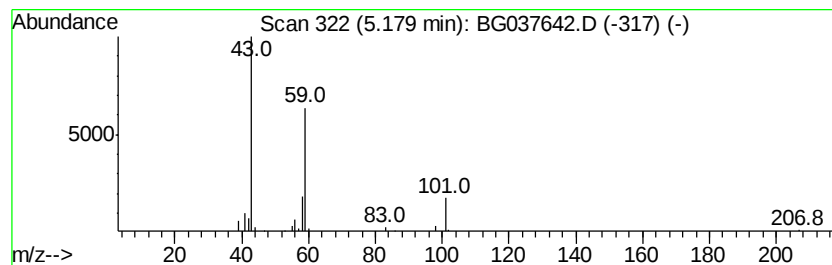
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.18	3.47 ng	52336	1,4-Dichlorobenzene-d4	8.10

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50
2			2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	28
3			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	10
4			Acetone	58	C3H6O	000067-64-1	9
5			Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9



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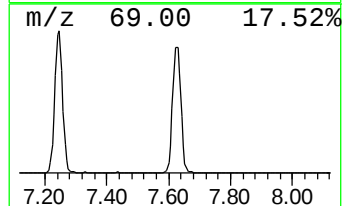
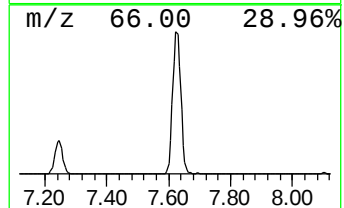
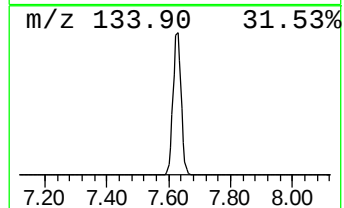
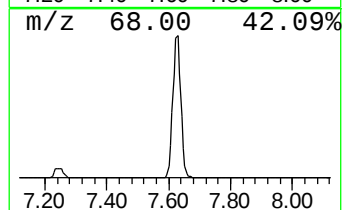
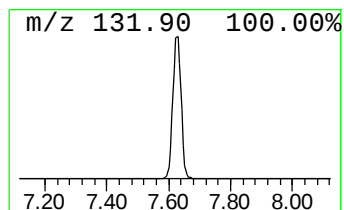
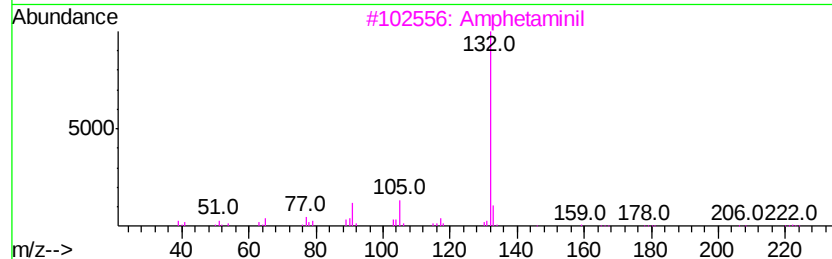
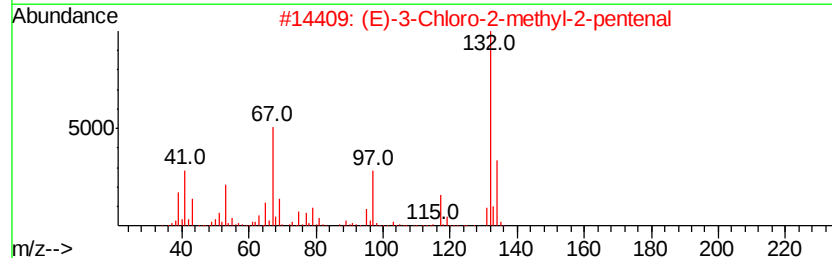
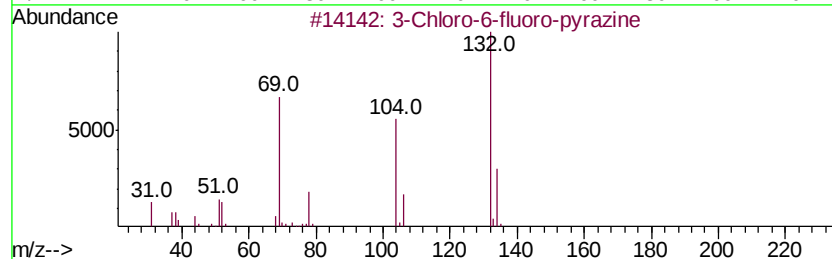
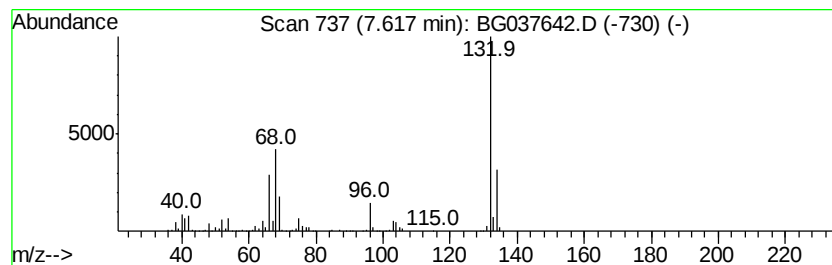
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Peak Number 2 unknown7.62 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.62	85.46 ng	1288110	1,4-Dichlorobenzene-d4	8.10

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	17
3		Amphetaminil	250	C17H18N2	017590-01-1	10
4		5-Aminoindole	132	C8H8N2	005192-03-0	10
5		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.18	3.5	ng	52336	1	8.10	301458	20.0
unknown7.62	7.62	85.5	ng	1288110	1	8.10	301458	20.0