

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG110218\
 Data File : BG037762.D
 Acq On : 2 Nov 2018 19:58
 Operator : JU/SJ
 Sample : PB114425BL
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB114425BL

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	5.181	315	322	335	rBV	49316	87756	1.93%	0.365%
2	5.639	393	400	412	rBV	1124684	1890438	41.51%	7.857%
3	7.243	665	673	685	rBV	1169379	2134706	46.87%	8.872%
4	7.625	727	738	752	rBV	1130758	2020456	44.36%	8.398%
5	8.095	811	818	827	rBV	217469	389507	8.55%	1.619%
6	8.424	865	874	885	rBV	832863	1540504	33.83%	6.403%
7	9.276	1010	1019	1031	rBV	716894	1325770	29.11%	5.510%
8	10.921	1291	1299	1309	rBV	325399	613151	13.46%	2.548%
9	13.354	1705	1713	1723	rBV	2100802	3353548	73.64%	13.938%
10	14.735	1941	1948	1955	rBV2	540284	900891	19.78%	3.744%
11	16.221	2193	2201	2224	rBV	1337625	2256766	49.55%	9.380%
12	17.478	2407	2415	2425	rBV2	678390	1068559	23.46%	4.441%
13	20.081	2851	2858	2871	rBV	3262852	4554278	100.00%	18.929%
14	21.762	3137	3144	3159	rBV2	572475	1102463	24.21%	4.582%
15	25.099	3700	3712	3735	rBV	220628	821158	18.03%	3.413%

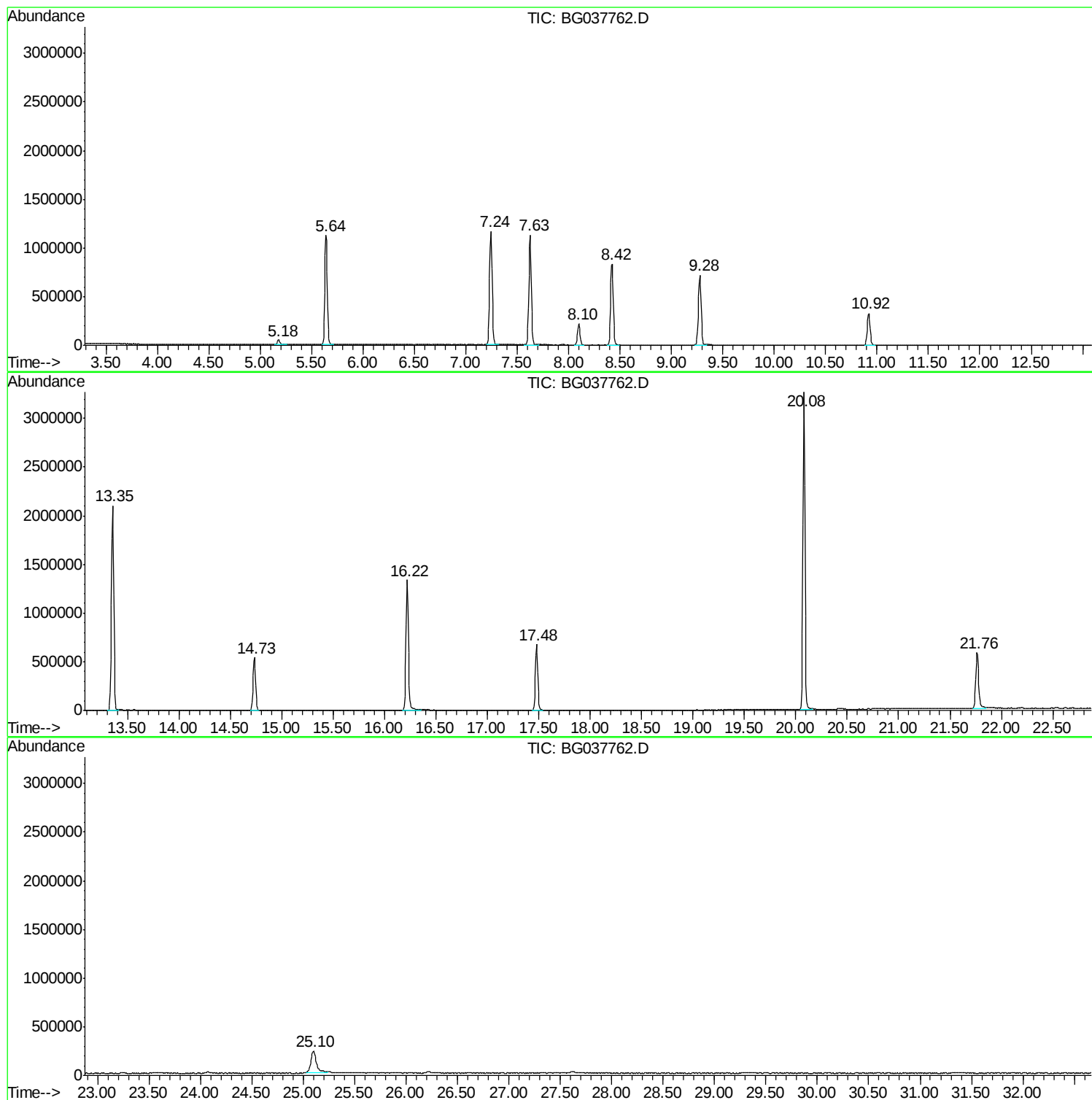
Sum of corrected areas: 24059951

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



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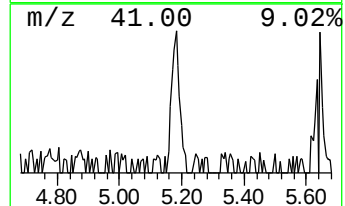
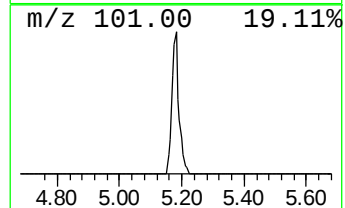
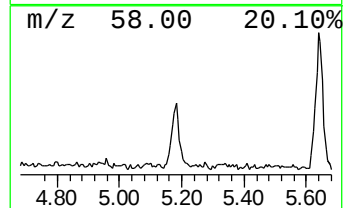
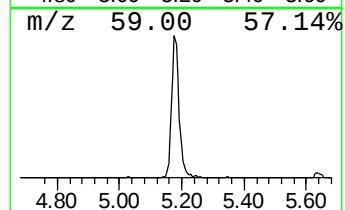
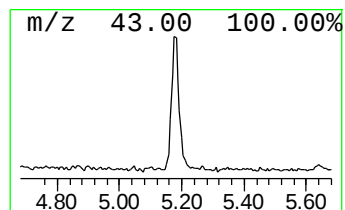
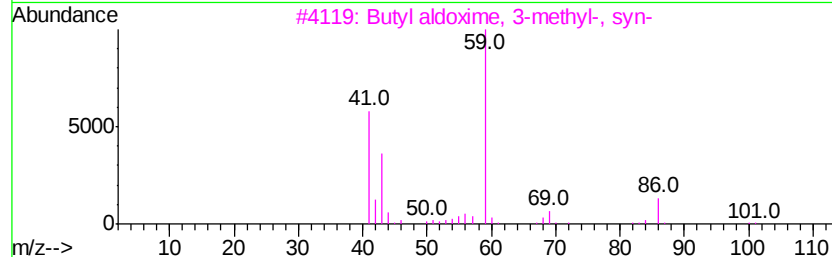
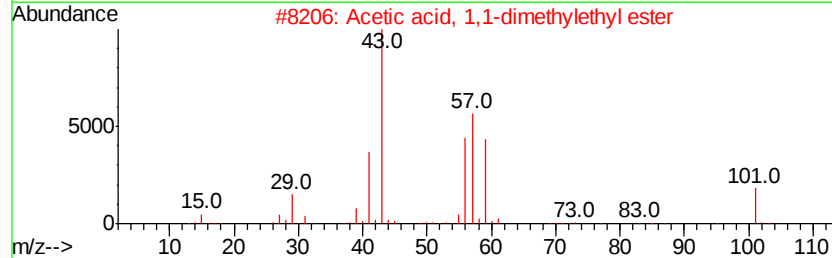
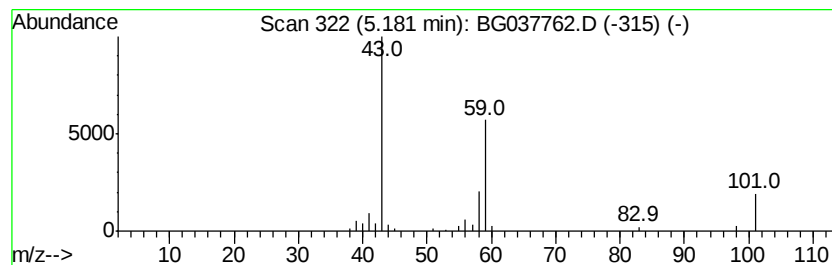
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.18	4.51 ng	87756	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	53
2			Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	38
3			Butyl aldoxime, 3-methyl-, syn-	101	C5H11NO	005780-40-5	25
4			Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	17
5			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9



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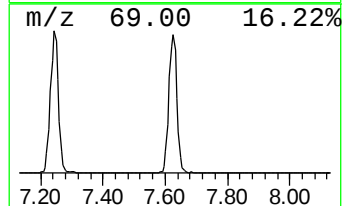
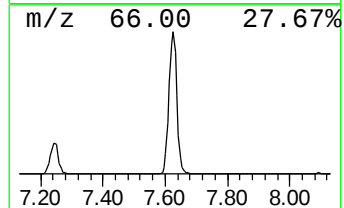
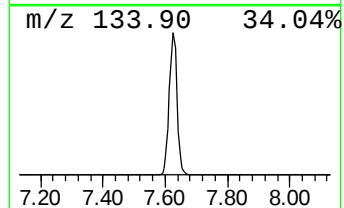
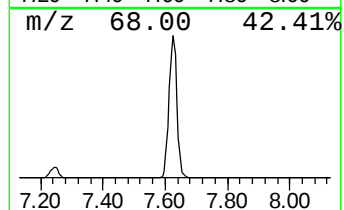
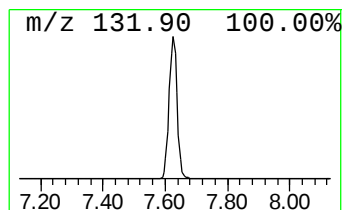
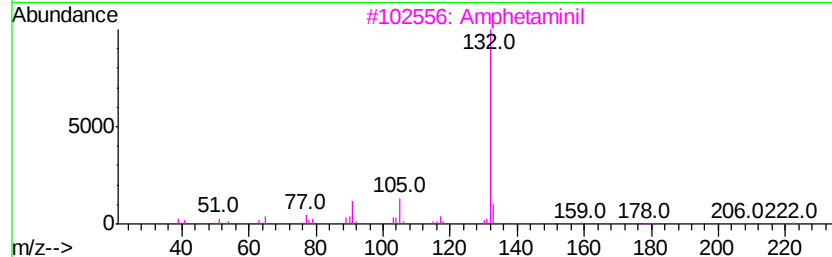
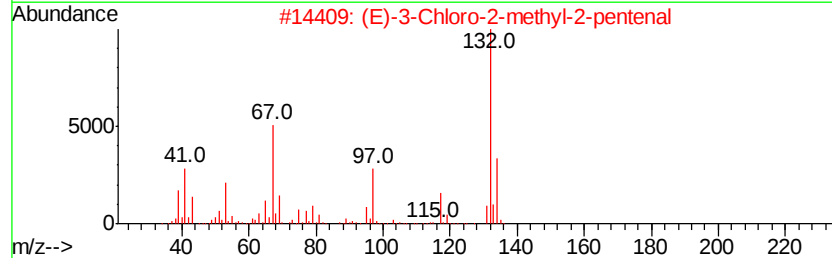
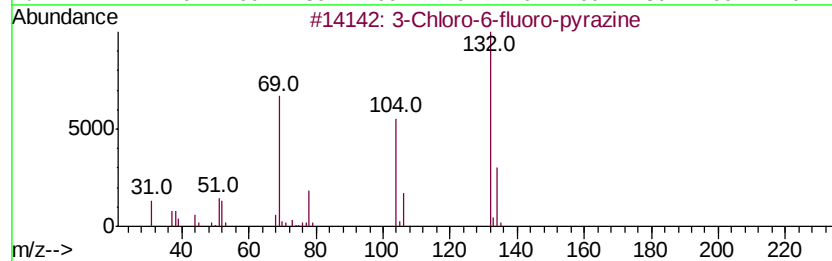
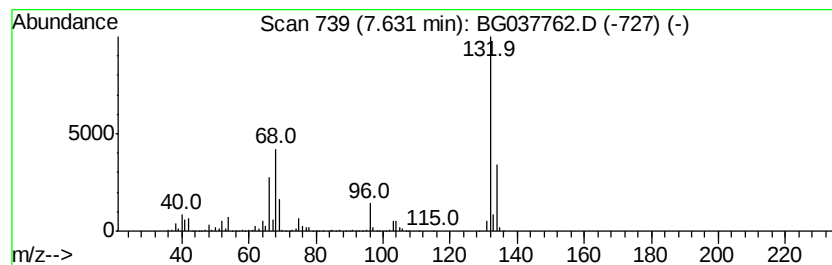
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Peak Number 2 unknown7.63 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.63	103.74 ng	2020460	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	16
3		Amphetaminil	250	C17H18N2	017590-01-1	12
4		Imidazole, 4,5-dicyano-1-methyl-	132	C6H4N4	019485-35-9	9
5		Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.18	4.5	ng	87756	1	8.10	389507	20.0
unknown7.63	7.63	103.7	ng	2020460	1	8.10	389507	20.0