

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG102318\
 Data File : BG037479.D
 Acq On : 23 Oct 2018 16:51
 Operator : JU/SJ
 Sample : PB114087BL
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB114087BL

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	5.186	317	323	331	rBV	30661	52576	1.67%	0.309%
2	5.644	393	401	409	rBV	732303	1190685	37.85%	6.996%
3	7.248	666	674	686	rBV	757087	1323575	42.07%	7.777%
4	7.630	732	739	753	rBV	707470	1229220	39.07%	7.223%
5	8.106	813	820	829	rVB	203422	350761	11.15%	2.061%
6	8.429	866	875	885	rVB	511796	922103	29.31%	5.418%
7	9.281	1011	1020	1031	rBV	416113	777117	24.70%	4.566%
8	10.926	1293	1300	1310	rVB	285626	523430	16.64%	3.076%
9	13.365	1707	1715	1724	rBV	1232329	1966966	62.52%	11.558%
10	14.740	1942	1949	1959	rBV2	495210	812843	25.84%	4.776%
11	16.226	2195	2202	2213	rBV	873570	1375507	43.72%	8.083%
12	17.483	2409	2416	2429	rBV2	668115	1103898	35.09%	6.487%
13	20.086	2853	2859	2870	rBV	2366017	3145990	100.00%	18.486%
14	21.773	3138	3146	3158	rBV	664574	1157041	36.78%	6.799%
15	25.110	3700	3714	3726	rBV2	328775	1086614	34.54%	6.385%

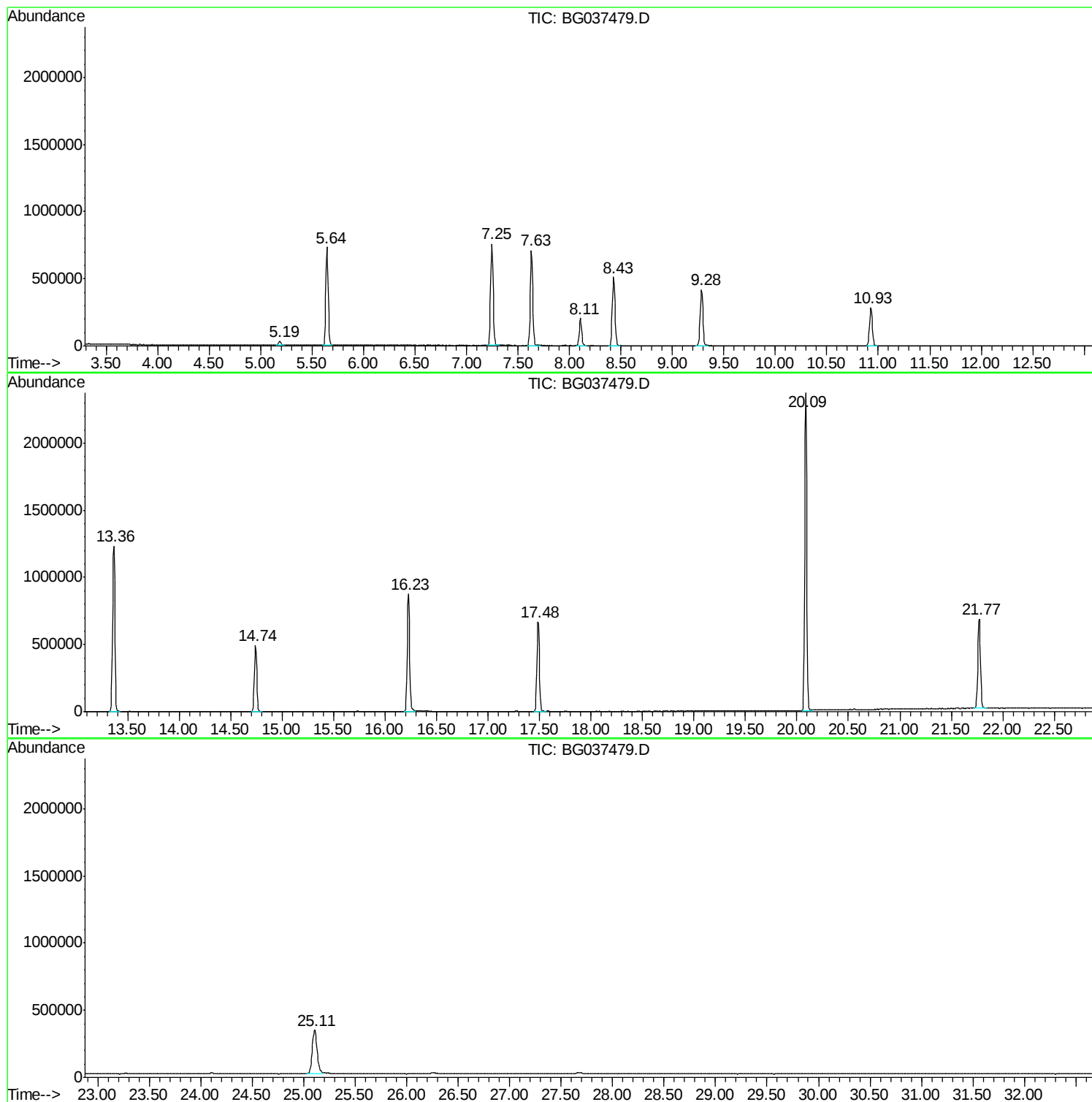
Sum of corrected areas: 17018326

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Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG101818.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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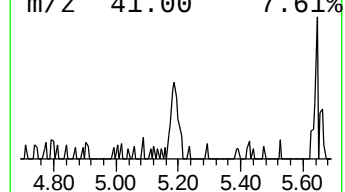
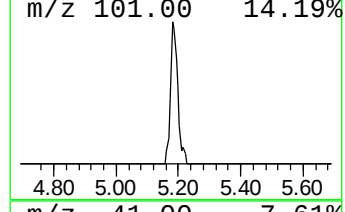
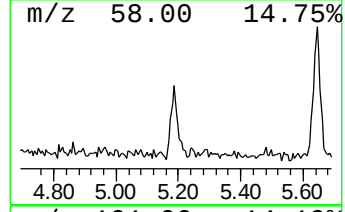
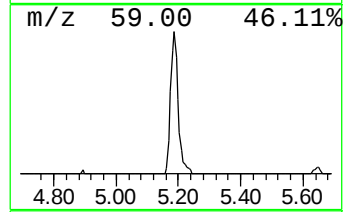
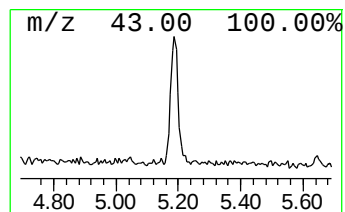
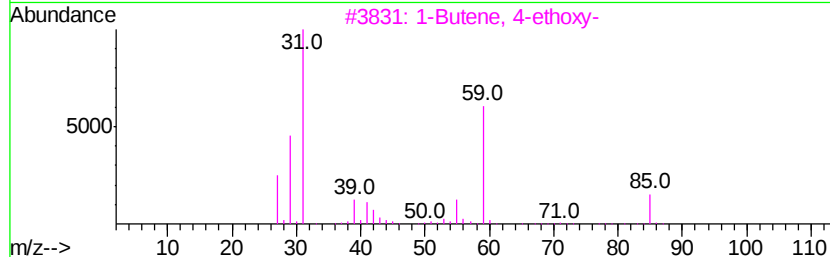
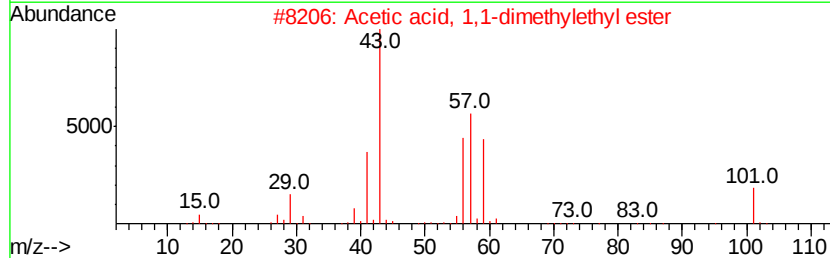
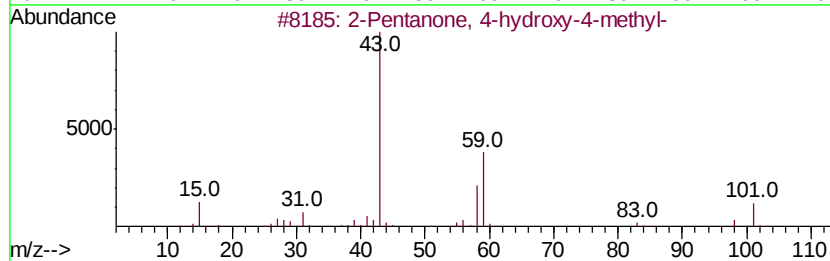
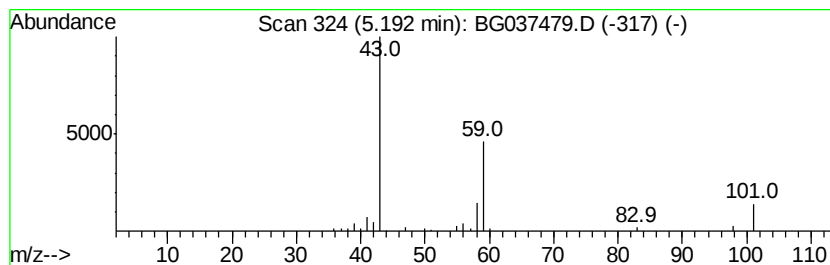
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.19	3.00 ng	52576	1,4-Dichlorobenzene-d4	8.11

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	83		
2	Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	36		
3	1-Butene, 4-ethoxy-	100	C6H12O	044611-46-3	17		
4	Diacetamide	101	C4H7NO2	000625-77-4	9		
5	5-Aminovaleric acid	117	C5H11NO2	000660-88-8	9		



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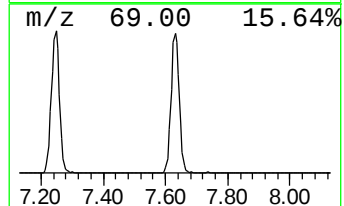
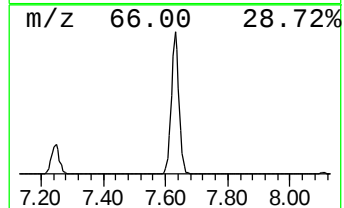
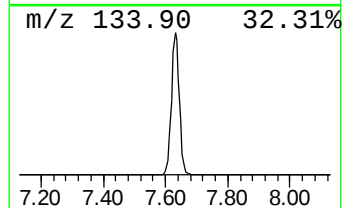
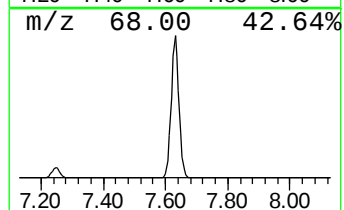
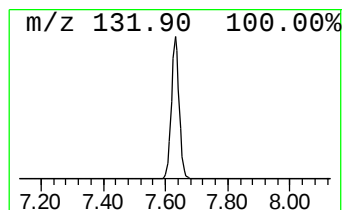
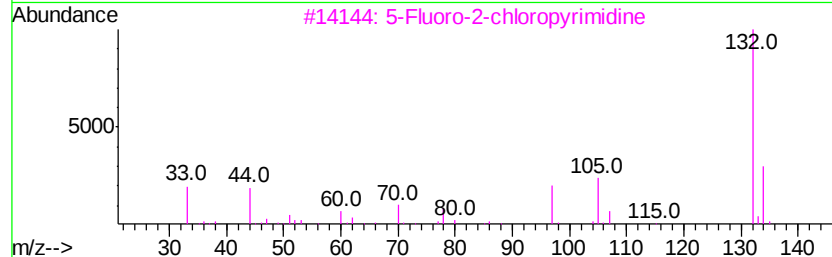
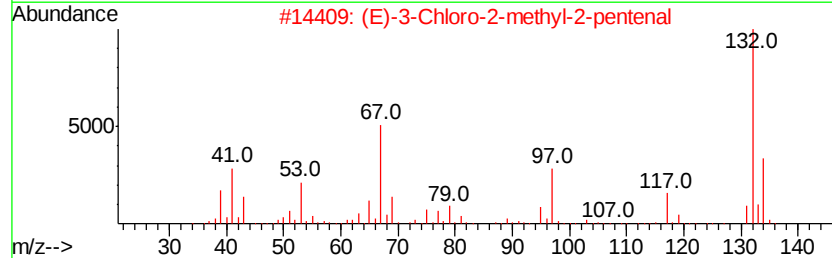
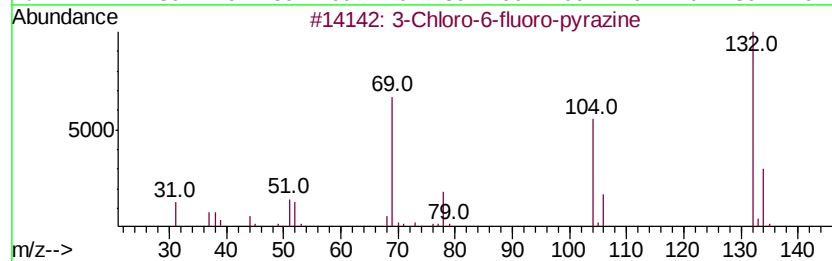
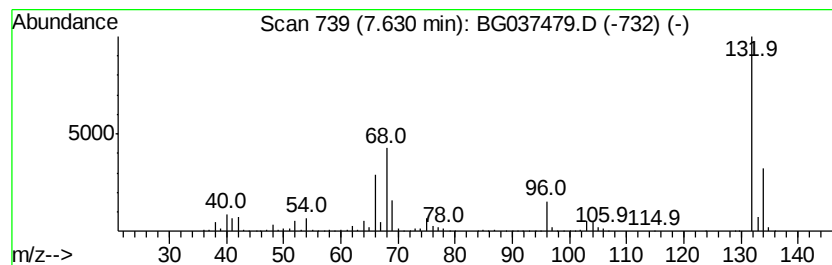
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Peak Number 2 unknown7.63 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.63	70.09 ng	1229220	1,4-Dichlorobenzene-d4	8.11

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		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	12
4		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
5		2H-Cyclopenta[d]pyridazine, 2-me...	132	C8H8N2	022291-85-6	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.19	3.0	ng	52576	1	8.11	350761	20.0
unknown7.63	7.63	70.1	ng	1229220	1	8.11	350761	20.0