

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG012519\
 Data File : BG039279.D
 Acq On : 25 Jan 2019 15:23
 Operator : JU/SJ
 Sample : PB116607BL
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB116607BL

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-BG010919.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	5.091	336	341	348	rBV	18224	30757	1.69%	0.414%
2	5.561	414	421	435	rVB	290488	480159	26.44%	6.455%
3	7.171	685	695	707	rBV	288479	513633	28.28%	6.905%
4	7.535	749	757	766	rBV	287875	495398	27.27%	6.660%
5	8.005	831	837	843	rBV	56732	96353	5.30%	1.295%
6	8.328	884	892	900	rBB	216605	379253	20.88%	5.099%
7	9.186	1030	1038	1049	rBV	196851	366524	20.18%	4.928%
8	10.825	1309	1317	1334	rBB	75018	144182	7.94%	1.938%
9	13.264	1725	1732	1742	rBV	699224	1104561	60.81%	14.850%
10	14.650	1960	1968	1979	rBV2	127952	222533	12.25%	2.992%
11	16.143	2215	2222	2237	rBV2	416379	677361	37.29%	9.107%
12	17.394	2429	2435	2447	rBV	190384	300112	16.52%	4.035%
13	20.003	2872	2879	2892	rBV	1412208	1816327	100.00%	24.419%
14	21.666	3155	3162	3171	rBV2	239328	412314	22.70%	5.543%
15	24.926	3707	3717	3728	rVB2	136535	398644	21.95%	5.359%

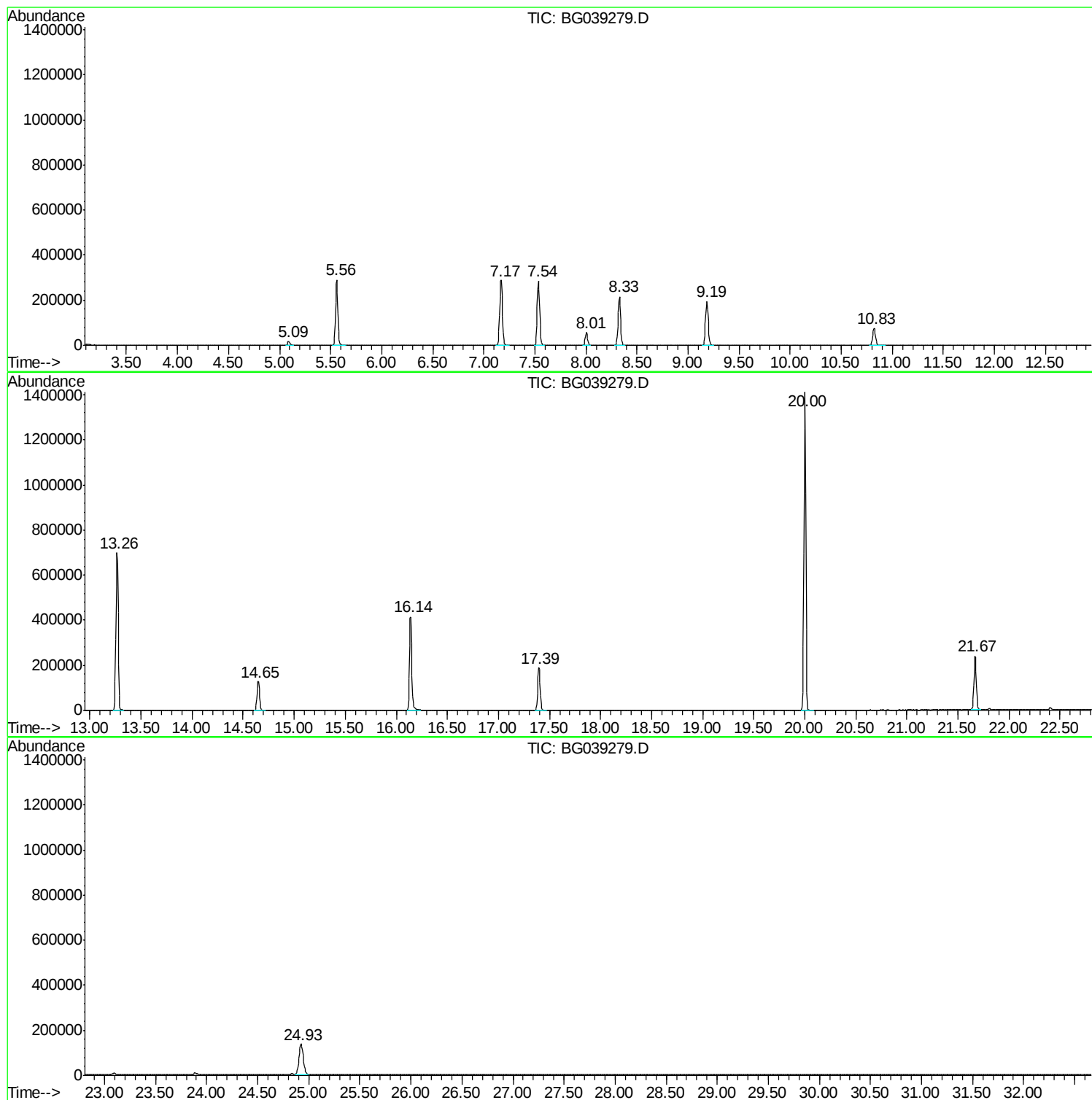
Sum of corrected areas: 7438111

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Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG010919.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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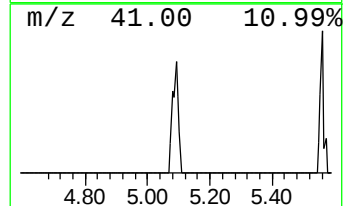
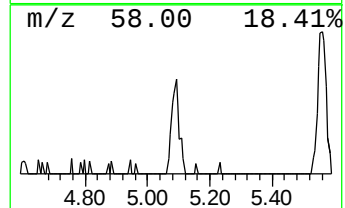
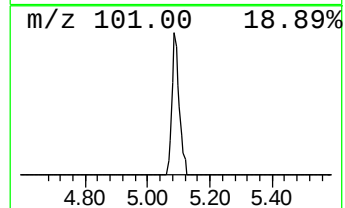
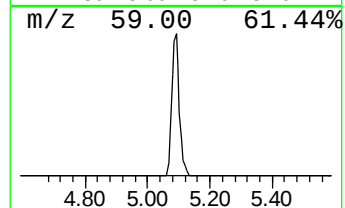
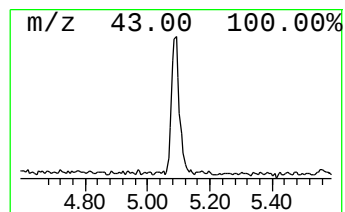
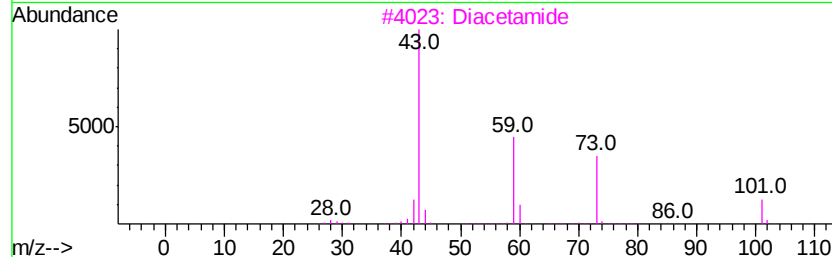
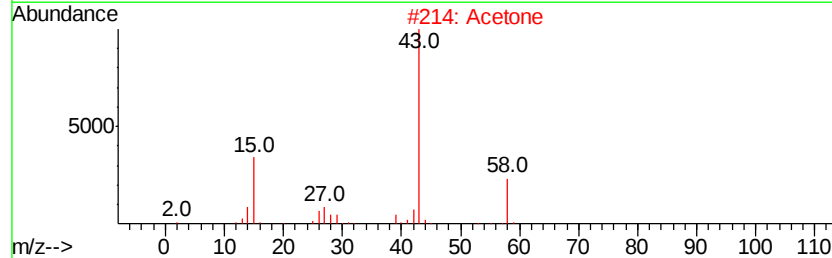
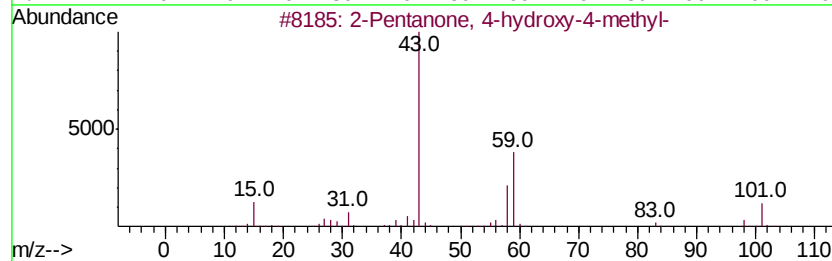
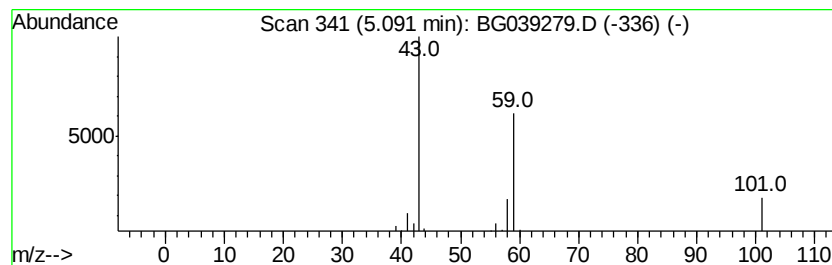
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.09	6.38 ng	30757	1,4-Dichlorobenzene-d4	8.01

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	78		
2	Acetone	58	C3H6O	000067-64-1	9		
3	Diacetamide	101	C4H7NO2	000625-77-4	9		
4	Propane, 2-methyl-2-(1-methyleth...	116	C7H16O	017348-59-3	9		
5	1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	9		



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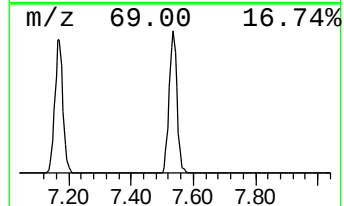
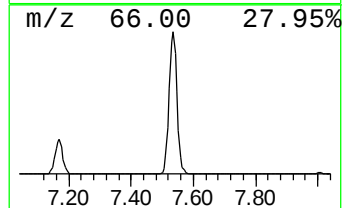
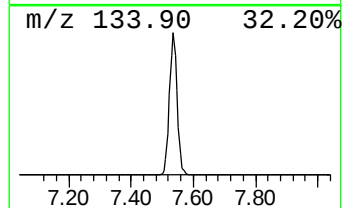
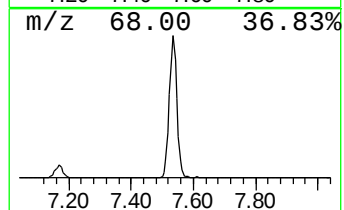
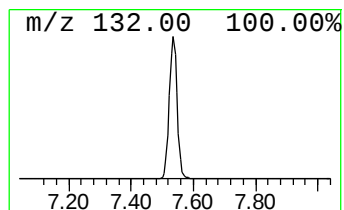
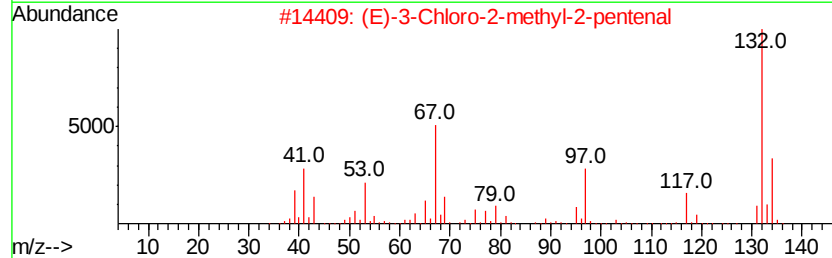
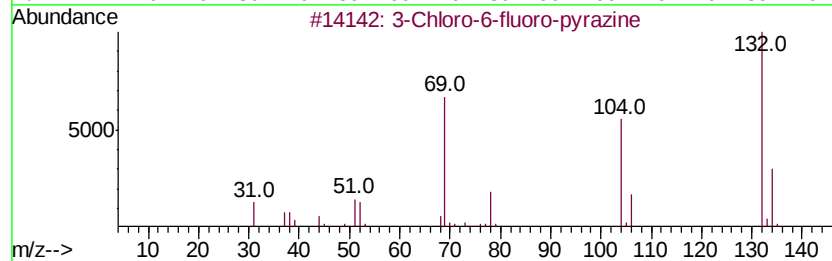
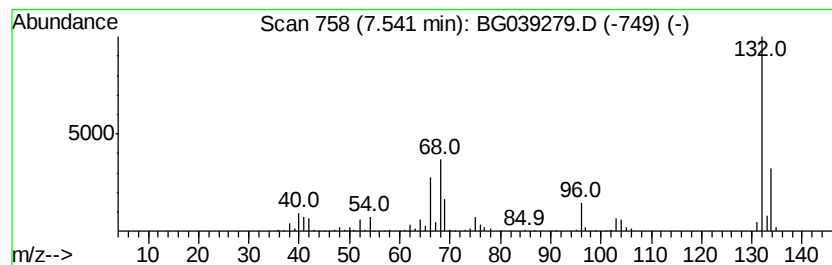
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Peak Number 2 unknown7.54 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.54	102.83 ng	495398	1,4-Dichlorobenzene-d4	8.01

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		5-Aminoindole	132	C8H8N2	005192-03-0	12
4		(5-Methyl-2-pyridyl)acetonitrile	132	C8H8N2	1000241-93-9	10
5		1,2,4-Trifluorobenzene	132	C6H3F3	000367-23-7	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.09	6.4	ng	30757	1	8.01	96353	20.0
unknown7.54	7.54	102.8	ng	495398	1	8.01	96353	20.0