

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG022619\
 Data File : BG039628.D
 Acq On : 27 Feb 2019 9:54
 Operator : JU/SJ
 Sample : PB117396BL
 Misc :
 ALS Vial : 31 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB117396BL

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-BG012919.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	5.246	361	367	375	rVB	75680	117823	3.25%	0.657%
2	5.710	439	446	455	rBV	668710	1065956	29.41%	5.943%
3	7.313	710	719	732	rBV	677869	1178350	32.51%	6.570%
4	7.695	775	784	793	rBV	635427	1107486	30.55%	6.175%
5	8.165	854	864	871	rVB	195429	328585	9.06%	1.832%
6	8.488	910	919	930	rBV	684192	1188052	32.78%	6.624%
7	9.340	1056	1064	1073	rBV	580171	1036115	28.58%	5.777%
8	10.985	1335	1344	1354	rBV	268318	484341	13.36%	2.700%
9	13.411	1750	1757	1769	rBV	1612499	2581794	71.23%	14.395%
10	14.791	1984	1992	1999	rBV2	441184	732530	20.21%	4.084%
11	16.272	2237	2244	2268	rBV	698717	1165488	32.15%	6.498%
12	17.535	2452	2459	2467	rBV2	622733	964904	26.62%	5.380%
13	20.125	2893	2900	2909	rBV	2717368	3624807	100.00%	20.210%
14	20.760	3003	3008	3011	rBV	41382	48870	1.35%	0.272%
15	21.817	3181	3188	3199	rVB2	680131	1146370	31.63%	6.392%
16	24.155	3578	3586	3592	rBV5	17042	37365	1.03%	0.208%
17	25.195	3748	3763	3779	rBV2	356470	1126836	31.09%	6.283%

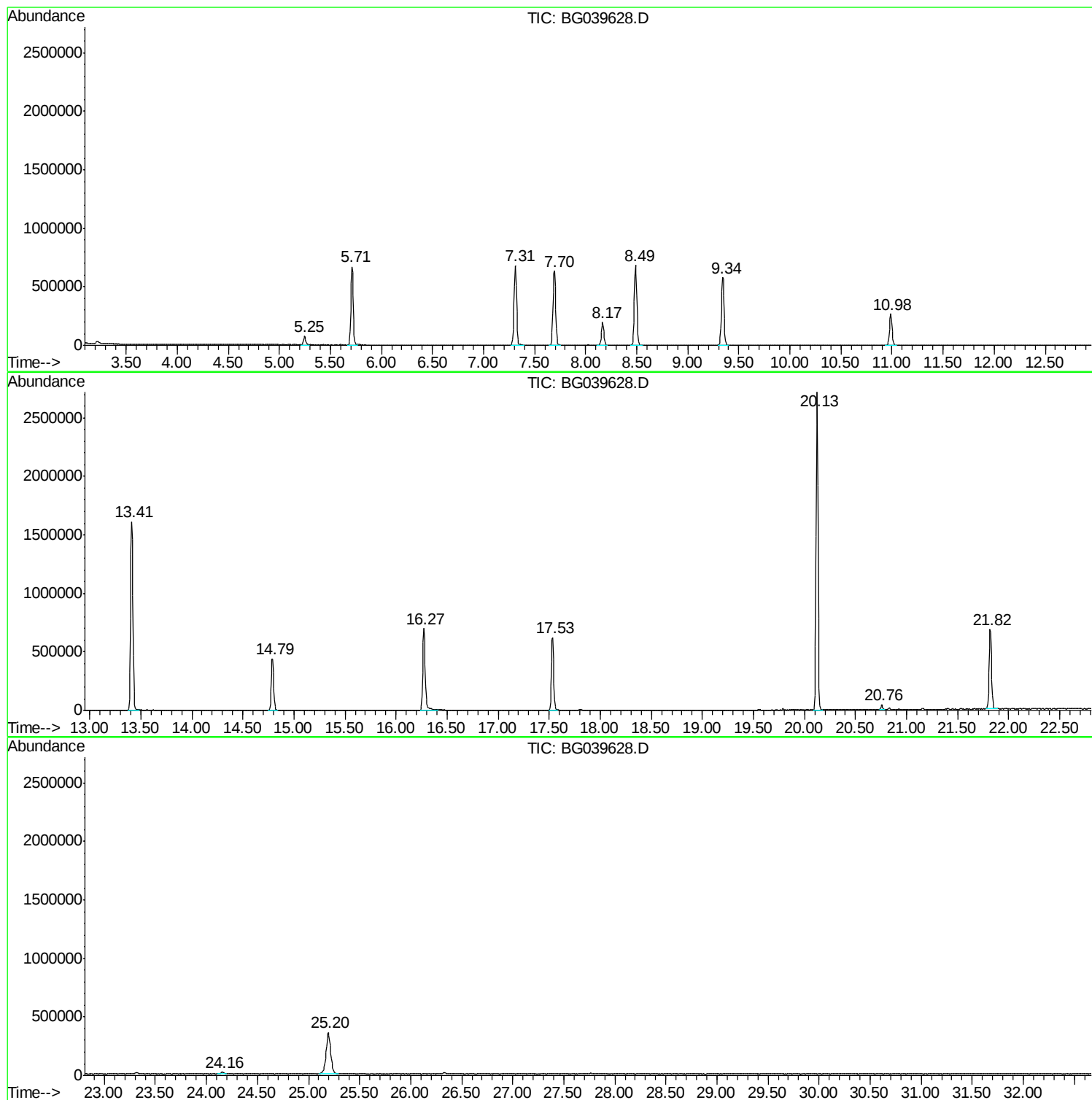
Sum of corrected areas: 17935672

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Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG012919.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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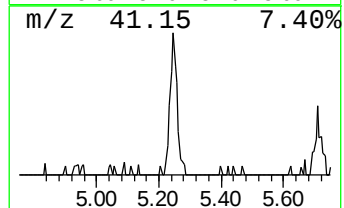
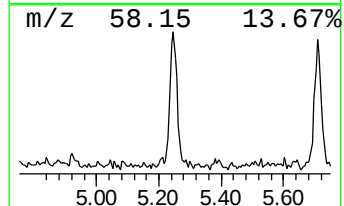
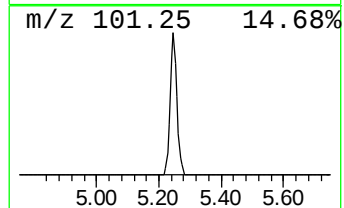
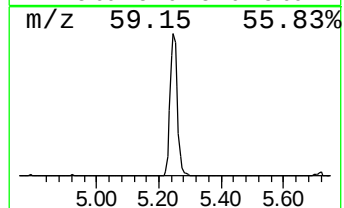
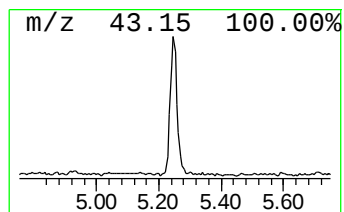
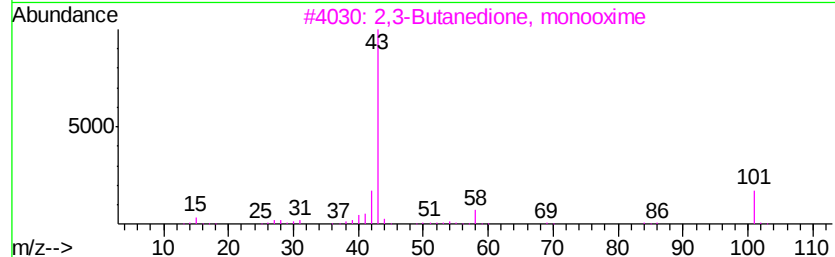
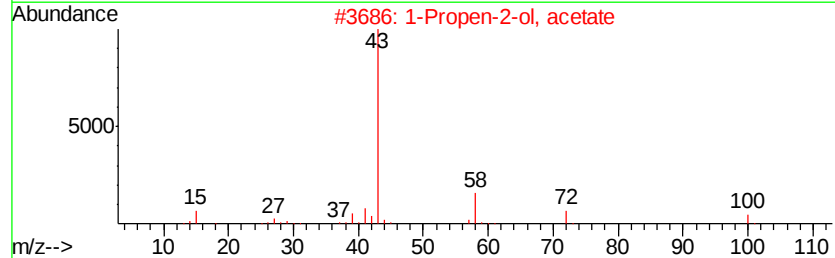
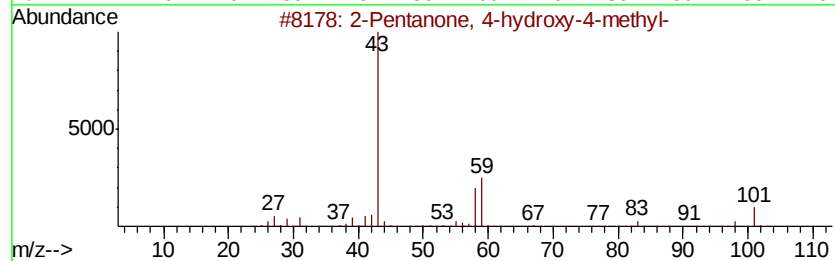
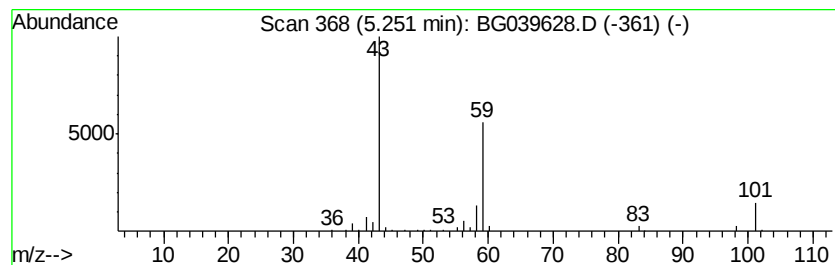
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.25	7.17 ng	117823	1,4-Dichlorobenzene-d4	8.17

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	64
2			1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	10
3			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	10
4			n-Propyl t-butyl ether	116	C7H16O	1000334-81-9	9
5			Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	9



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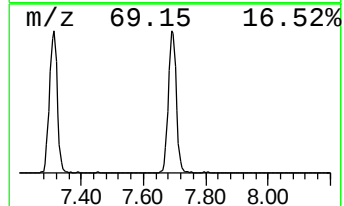
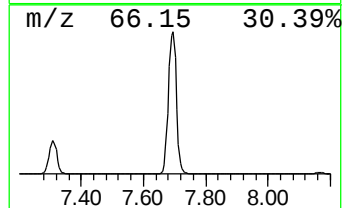
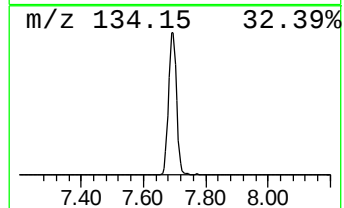
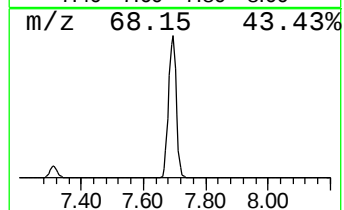
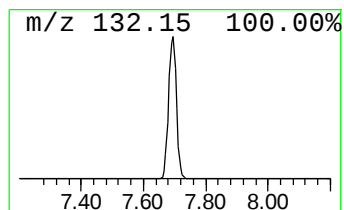
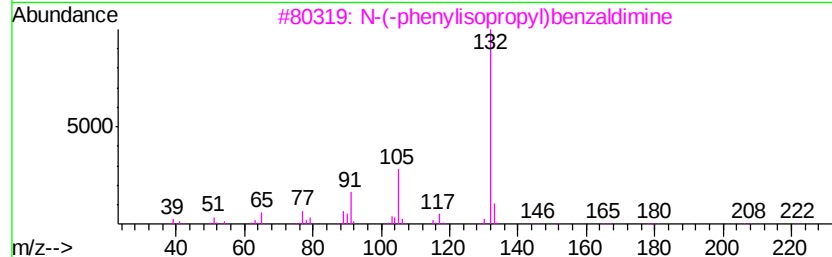
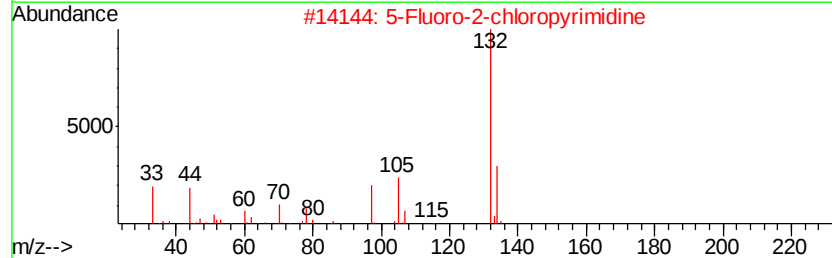
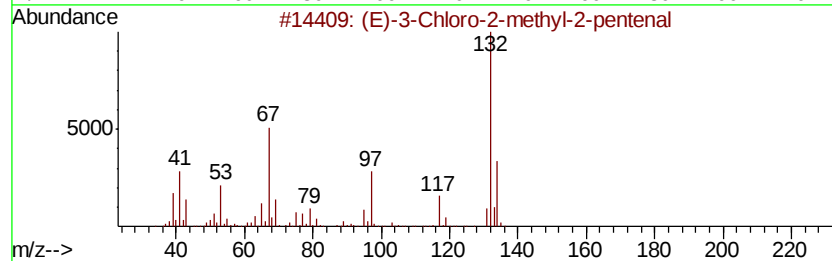
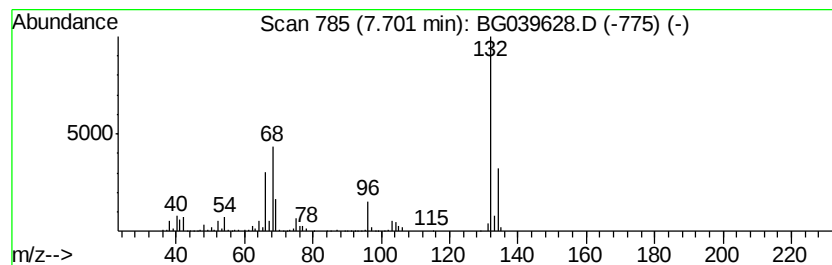
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Peak Number 2 unknown7.70 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.70	67.41 ng	1107490	1,4-Dichlorobenzene-d4	8.17

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	16
2		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	12
3		N-(-phenylisopropyl)benzalimine	223	C16H17N	1000379-01-3	10
4		Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9
5		Xenon	132	Xe	007440-63-3	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.25	7.2	ng	117823	1	8.17	328585	20.0
unknown7.70	7.70	67.4	ng	1107490	1	8.17	328585	20.0