

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG010919\
 Data File : BG039039.D
 Acq On : 10 Jan 2019 3:01
 Operator : JU/SJ
 Sample : PB116262BL
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB116262BL

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	5.111	339	344	350	rBV	11077	18396	1.18%	0.272%
2	5.581	417	424	434	rBV	233727	381403	24.49%	5.639%
3	7.185	690	697	708	rBV	238022	431937	27.73%	6.386%
4	7.555	752	760	769	rBV	227824	413206	26.53%	6.109%
5	8.025	834	840	850	rVB	62347	110667	7.11%	1.636%
6	8.348	887	895	904	rBV	170693	302590	19.43%	4.474%
7	9.206	1031	1041	1050	rBV	134624	253888	16.30%	3.754%
8	10.845	1311	1320	1331	rBV	90612	167890	10.78%	2.482%
9	13.289	1729	1736	1745	rBV	462159	734090	47.13%	10.853%
10	14.670	1964	1971	1981	rBV2	175418	290910	18.68%	4.301%
11	16.163	2219	2225	2243	rBV	441373	683777	43.90%	10.110%
12	17.420	2431	2439	2451	rBV	292296	444147	28.52%	6.567%
13	20.023	2876	2882	2889	rBV	1265133	1557506	100.00%	23.027%
14	21.697	3160	3167	3179	rBV	302861	499397	32.06%	7.384%
15	24.982	3715	3726	3737	rVB	161665	473875	30.43%	7.006%

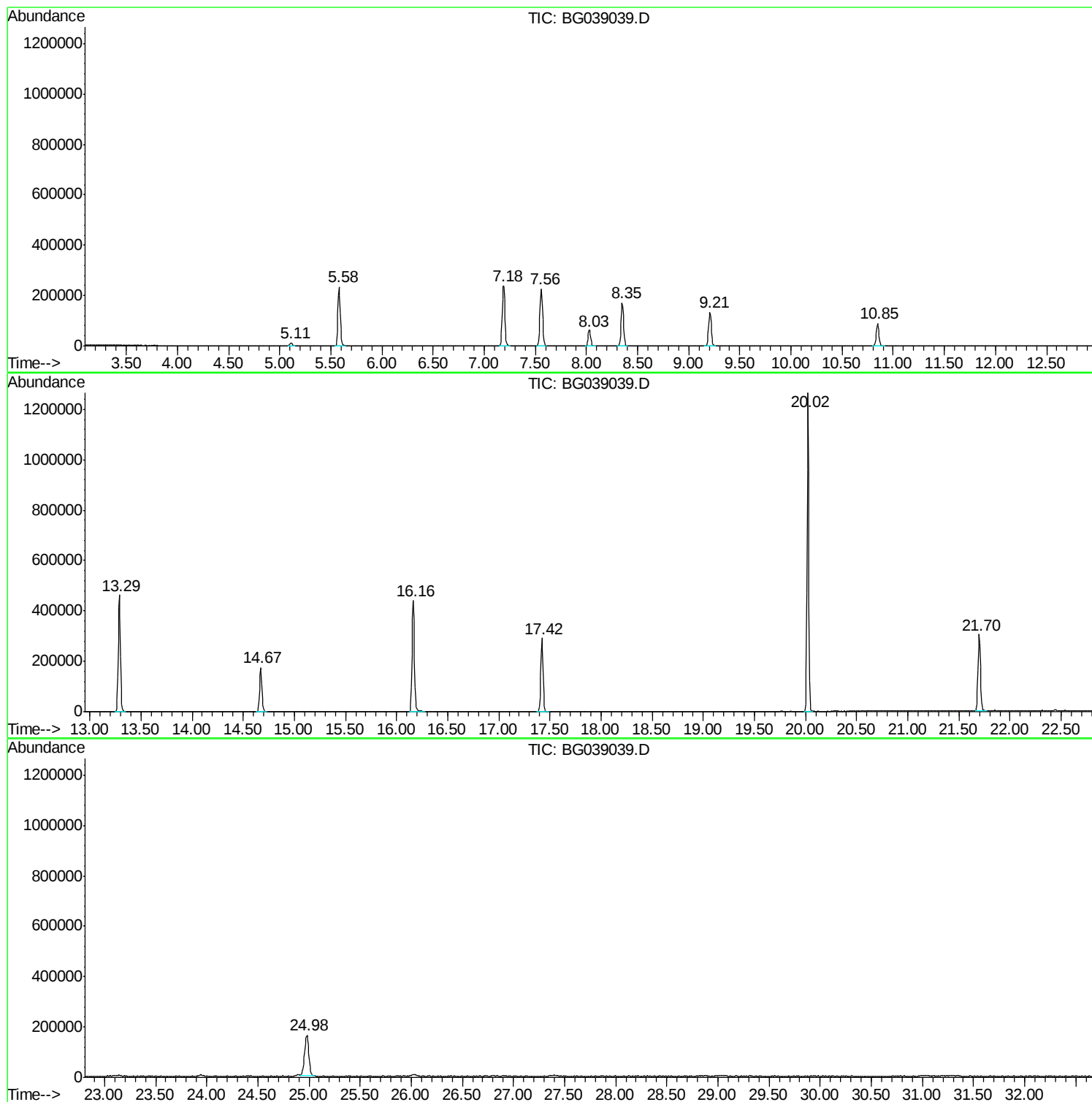
Sum of corrected areas: 6763679

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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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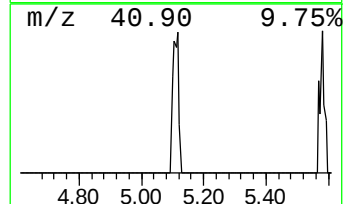
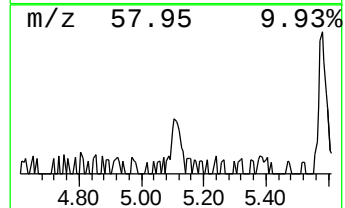
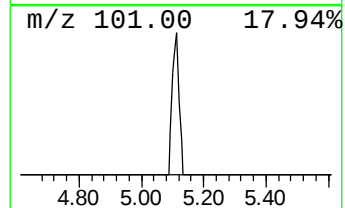
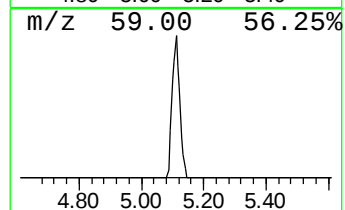
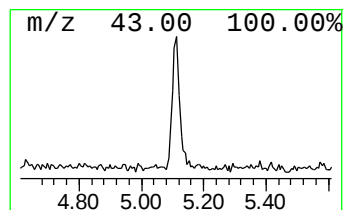
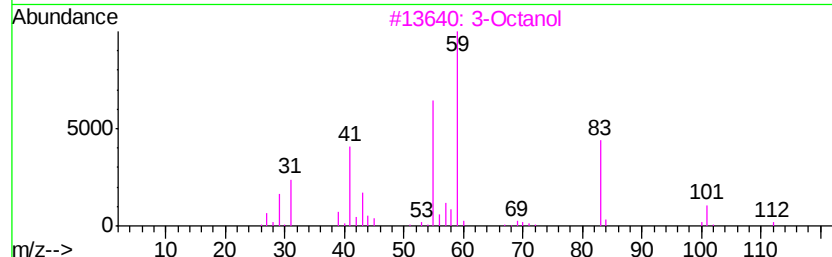
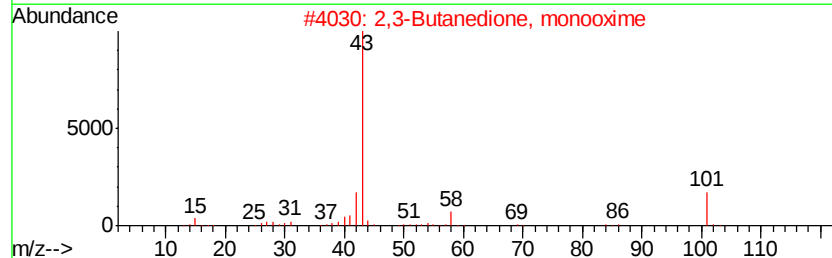
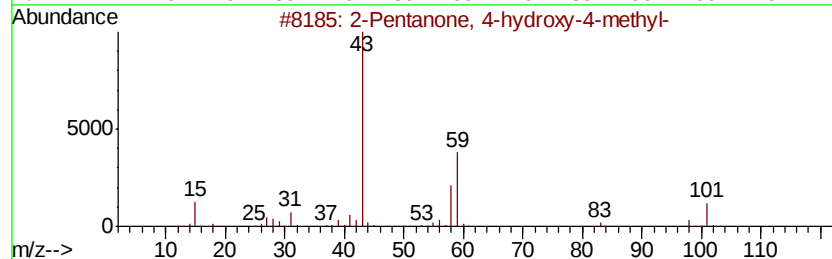
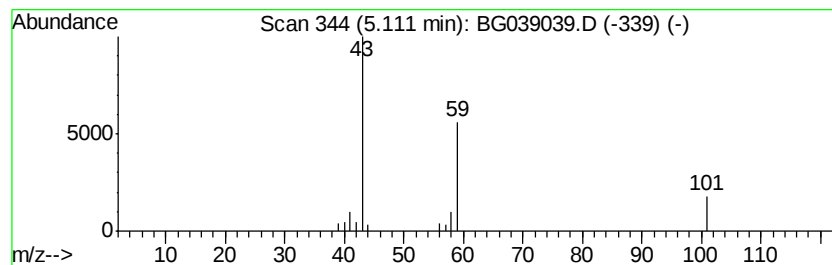
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TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.11	3.32 ng	18396	1,4-Dichlorobenzene-d4	8.03

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56
2	2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9
3	3-Octanol	130	C8H18O	000589-98-0	9
4	Diacetamide	101	C4H7NO2	000625-77-4	7
5	Butane	58	C4H10	000106-97-8	5



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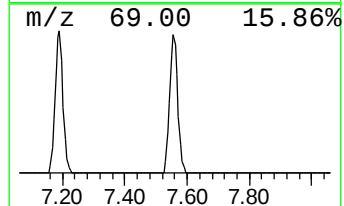
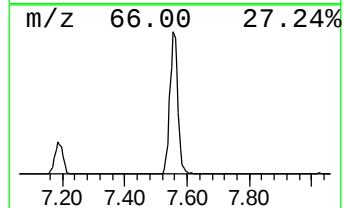
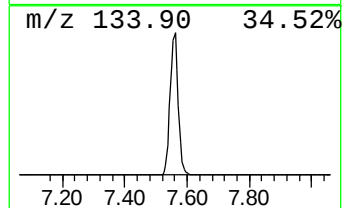
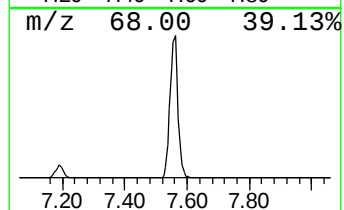
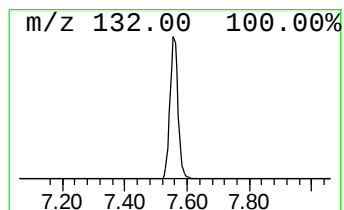
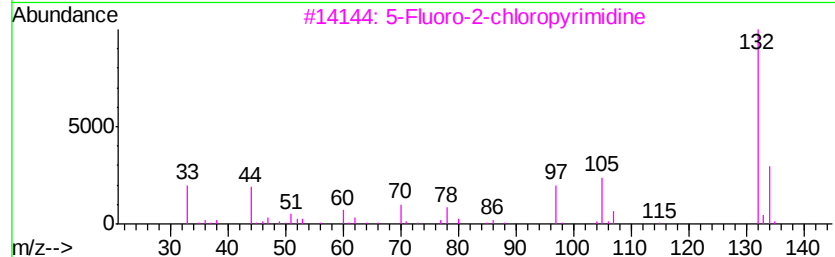
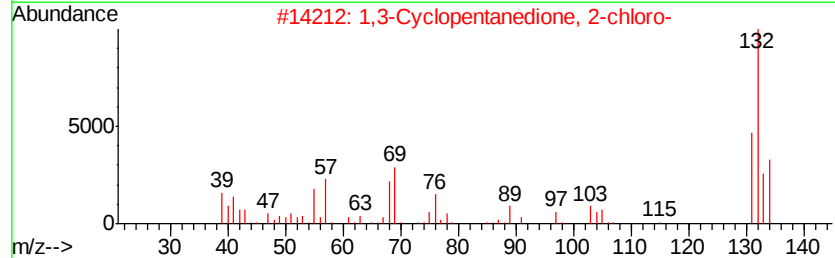
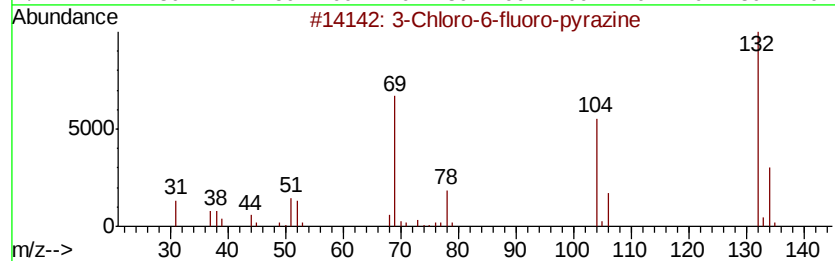
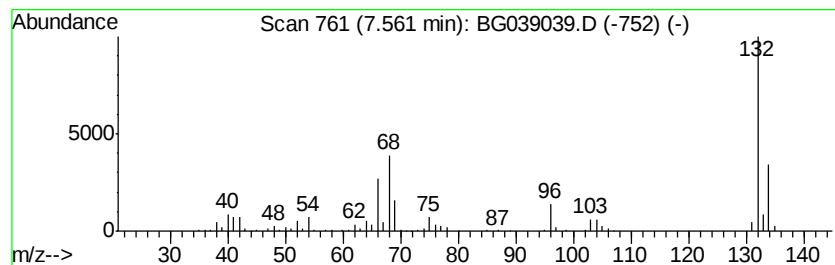
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Peak Number 2 unknown7.56 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.56	74.68 ng	413206	1,4-Dichlorobenzene-d4	8.03

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	35
2		1,3-Cyclopentanedione, 2-chloro-	132	C5H5ClO2	014203-19-1	32
3		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	17
4		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	17
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...	5.11	3.3	ng	18396	1	8.03	110667	20.0
unknown7.56	7.56	74.7	ng	413206	1	8.03	110667	20.0