

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG020119\
 Data File : BG039347.D
 Acq On : 1 Feb 2019 17:54
 Operator : JU/SJ
 Sample : PB116762BL
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB116762BL

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	5.258	364	369	380	rBV	57183	90659	2.22%	0.465%
2	5.716	438	447	456	rBV	1020622	1598686	39.23%	8.198%
3	7.314	711	719	732	rBV	1019661	1782950	43.75%	9.143%
4	7.702	776	785	796	rVB	968097	1648667	40.46%	8.455%
5	8.178	858	866	874	rBV	160357	276690	6.79%	1.419%
6	8.501	912	921	930	rBV	681174	1187540	29.14%	6.090%
7	9.352	1058	1066	1078	rBV	562809	1001357	24.57%	5.135%
8	10.997	1337	1346	1356	rVB	221199	404015	9.91%	2.072%
9	13.429	1752	1760	1769	rBV	1614586	2495566	61.24%	12.798%
10	14.804	1987	1994	2003	rVB2	356297	579370	14.22%	2.971%
11	16.284	2239	2246	2262	rBV	1179261	1824730	44.78%	9.358%
12	17.541	2454	2460	2468	rVB2	481292	739941	18.16%	3.795%
13	20.138	2896	2902	2908	rBV	3047267	4074974	100.00%	20.897%
14	21.836	3184	3191	3199	rVB2	540615	920150	22.58%	4.719%
15	25.219	3757	3767	3778	rVB	303884	874610	21.46%	4.485%

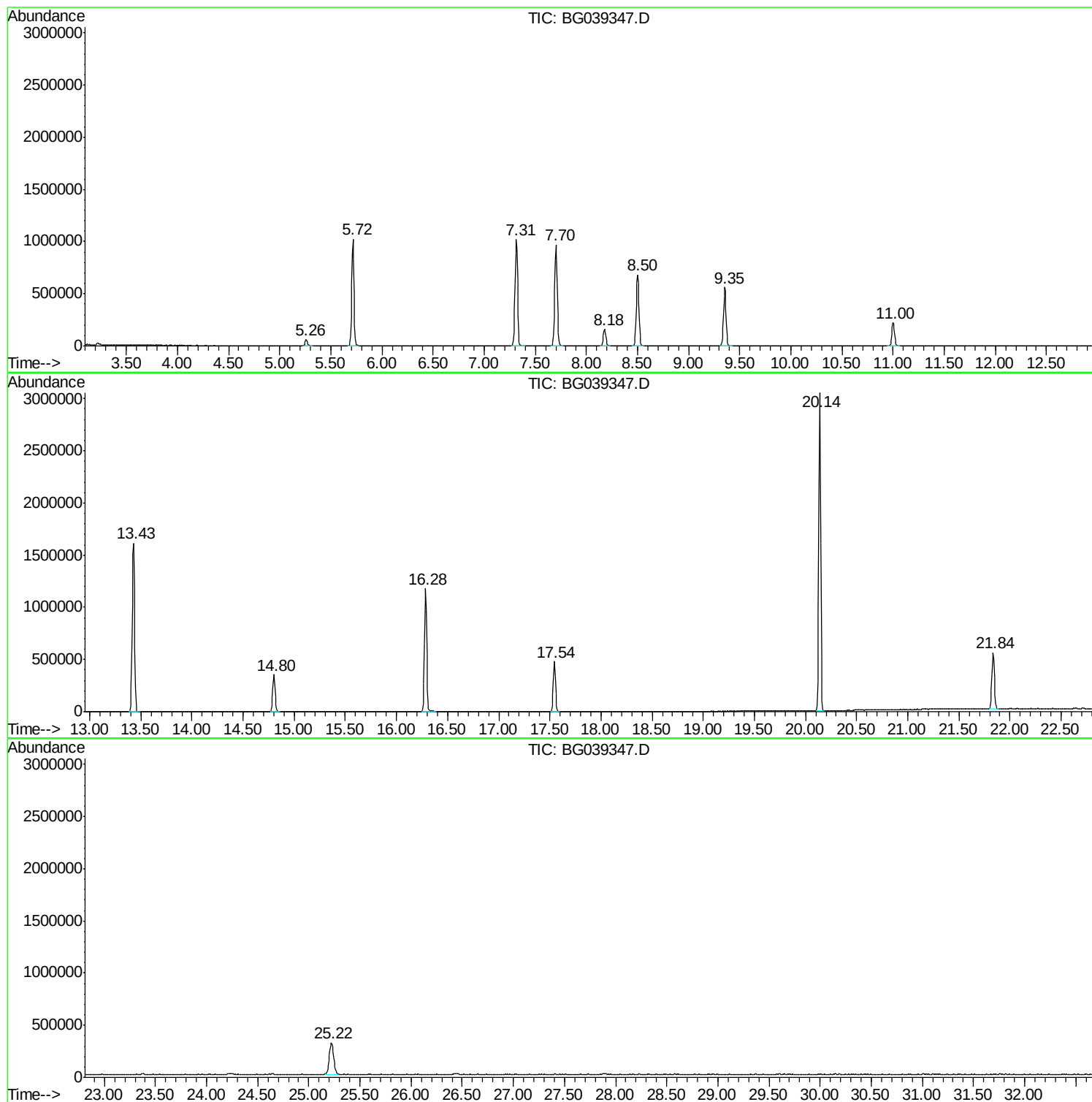
Sum of corrected areas: 19499905

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Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG012919.M
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TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P



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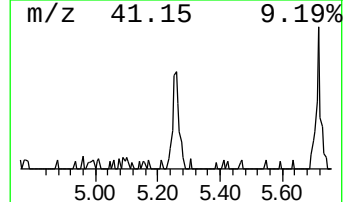
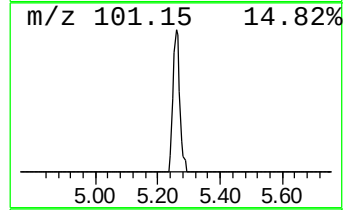
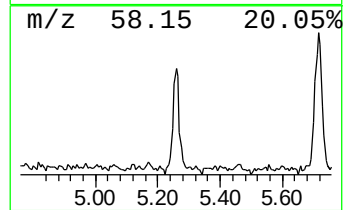
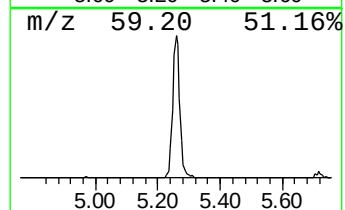
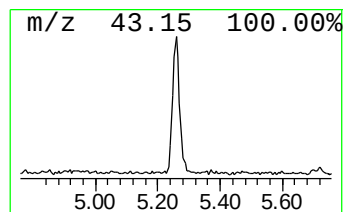
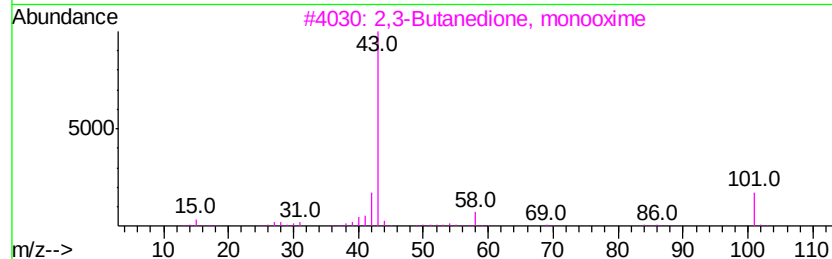
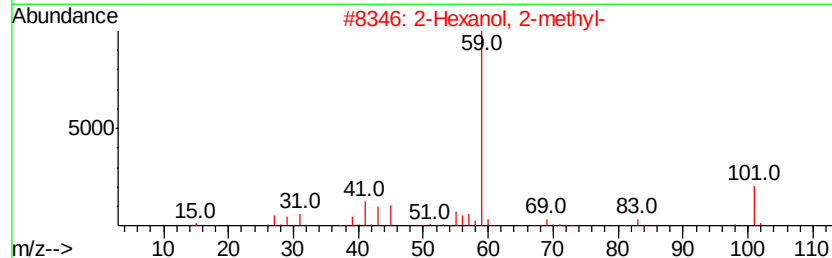
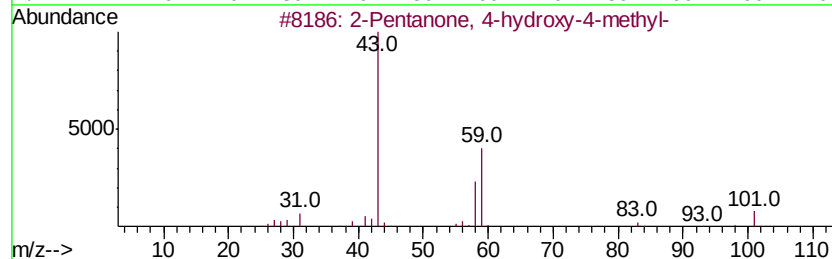
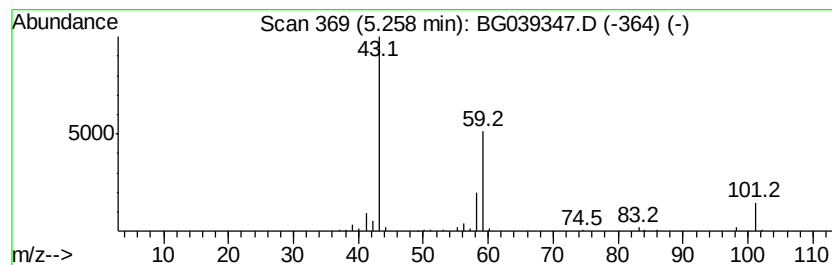
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.26	6.55 ng	90659	1,4-Dichlorobenzene-d4	8.18

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56
2		2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	25
3		2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	22
4		Butyl aldoxime, 3-methyl-, syn-	101	C5H11NO	005780-40-5	17
5		Acetone	58	C3H6O	000067-64-1	16



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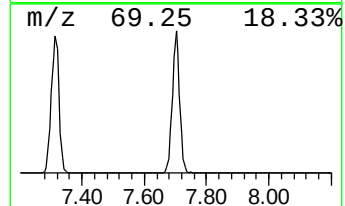
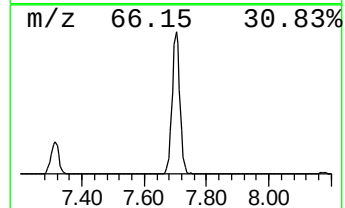
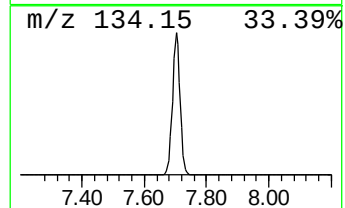
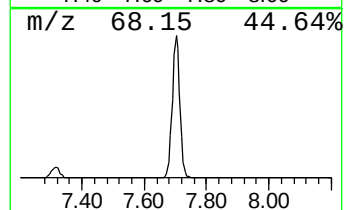
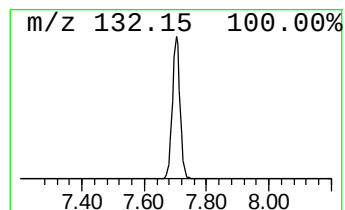
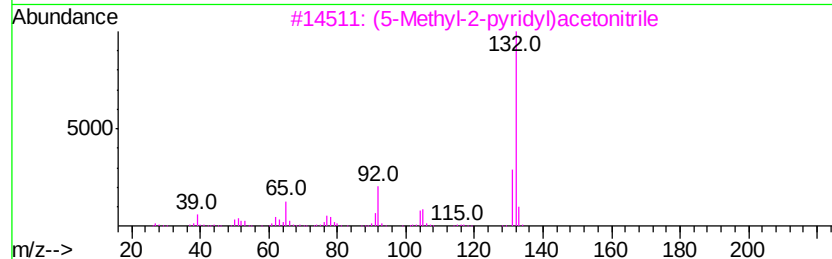
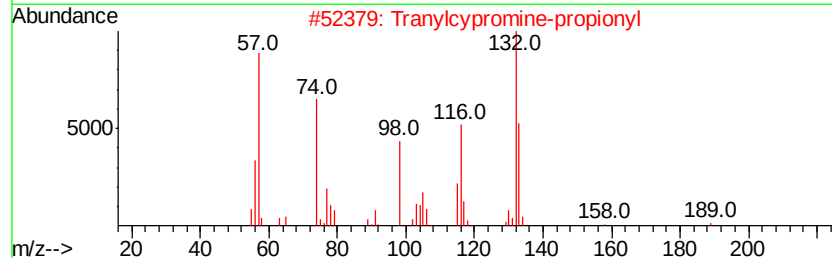
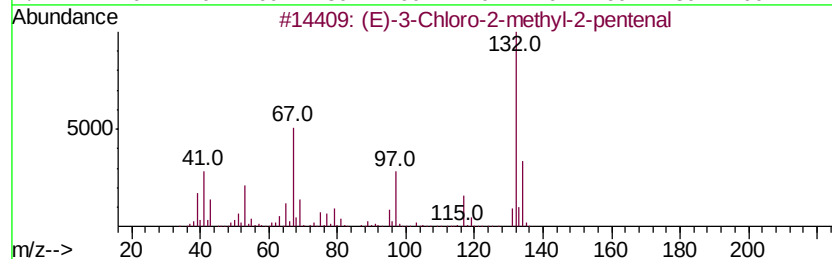
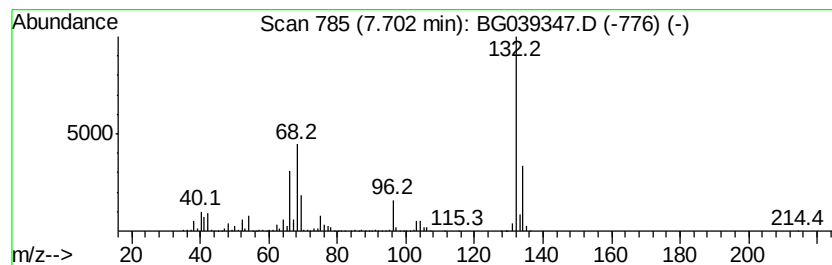
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Peak Number 2 unknown7.70 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.70	119.17 ng	1648670	1,4-Dichlorobenzene-d4	8.18

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	16
2		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	12
3		(5-Methyl-2-pyridyl)acetonitrile	132	C8H8N2	1000241-93-9	10
4		Imidazole, 4,5-dicyano-1-methyl-	132	C6H4N4	019485-35-9	9
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...	5.26	6.5	ng	90659	1	8.18	276690	20.0
unknown7.70	7.70	119.2	ng	1648670	1	8.18	276690	20.0