

Data Path : Z:\SVOASRV\HPCHEM1\BNA_G\DATA\BG101918\
 Data File : BG037447.D
 Acq On : 19 Oct 2018 10:45
 Operator : JU/SJ
 Sample : PB113692BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB113692BL

Integration Parameters: rteint.p
 Integrator: RTE
 Smoothing : ON
 Sampling : 1
 Start Thrs: 0.2
 Stop Thrs : 0
 Filtering: 5
 Min Area: 3 % of largest Peak
 Max Peaks: 100
 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	5.185	318	323	331	rBV	38252	63294	1.49%	0.291%
2	5.644	392	401	410	rBV	983673	1599693	37.54%	7.343%
3	7.248	666	674	688	rBV	1031838	1796032	42.15%	8.244%
4	7.630	730	739	748	rBV	937516	1660503	38.97%	7.622%
5	8.105	813	820	829	rVB	210027	365444	8.58%	1.677%
6	8.429	866	875	886	rBV	667160	1206428	28.31%	5.537%
7	9.281	1011	1020	1035	rBV	576345	1048836	24.61%	4.814%
8	10.926	1291	1300	1310	rVB	296362	555324	13.03%	2.549%
9	13.364	1707	1715	1724	rBV	1678826	2677352	62.83%	12.289%
10	14.739	1942	1949	1957	rBV2	527292	866271	20.33%	3.976%
11	16.225	2195	2202	2220	rBV4	1345304	2044215	47.97%	9.383%
12	17.483	2409	2416	2430	rVB2	734747	1184037	27.79%	5.435%
13	20.086	2853	2859	2871	rBV	3059740	4261098	100.00%	19.558%
14	21.766	3139	3145	3157	rVB2	710700	1250087	29.34%	5.738%
15	25.103	3702	3713	3734	rBV	374067	1208127	28.35%	5.545%

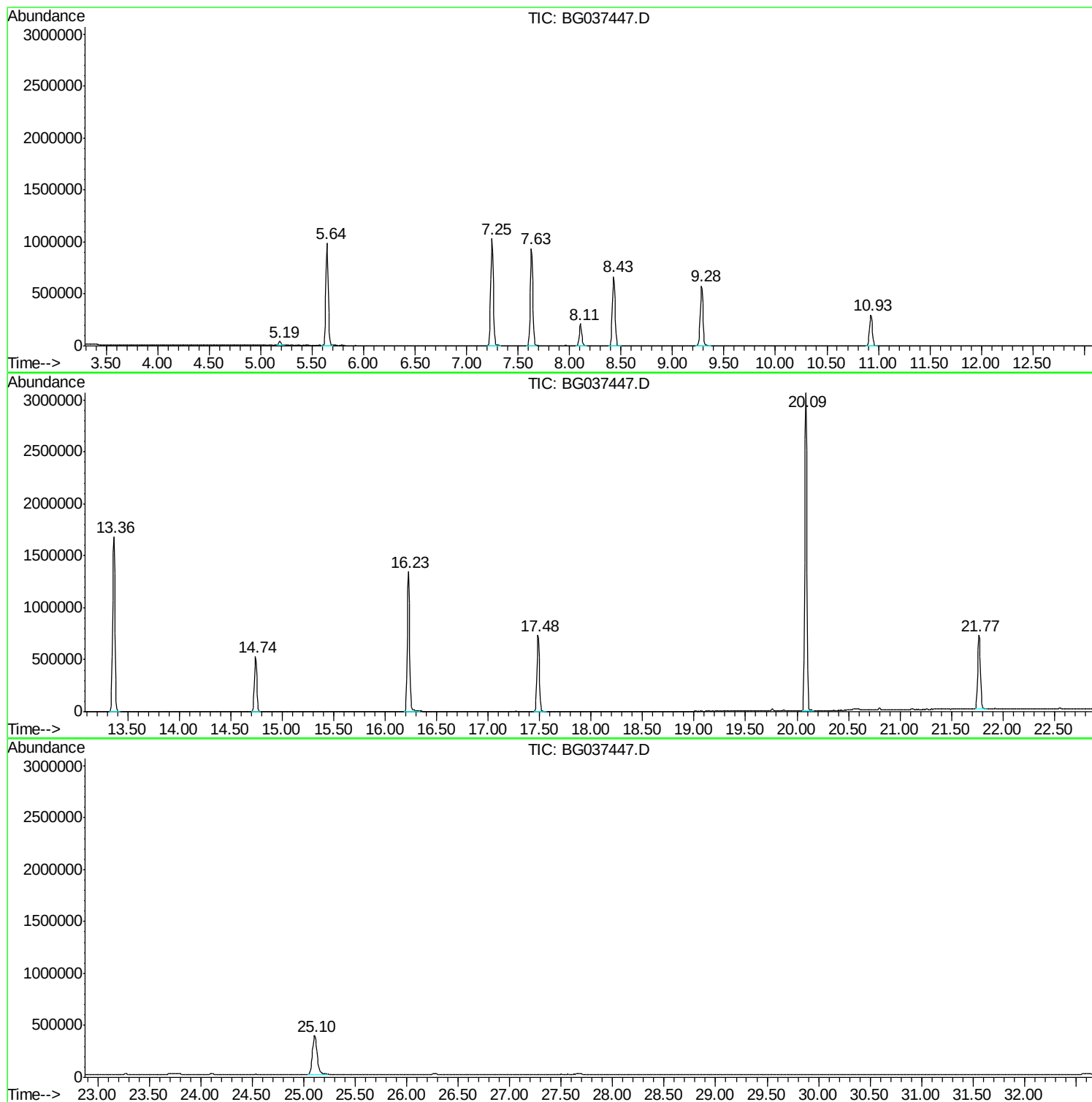
Sum of corrected areas: 21786741

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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG101818.M
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TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P



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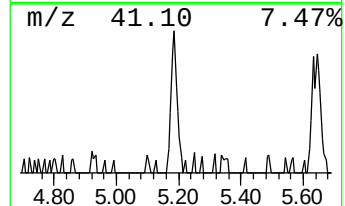
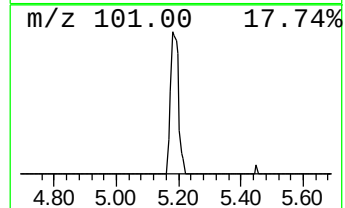
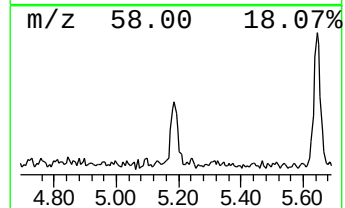
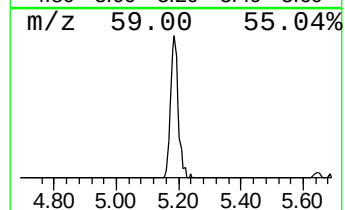
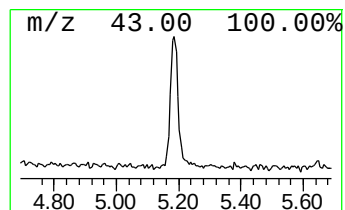
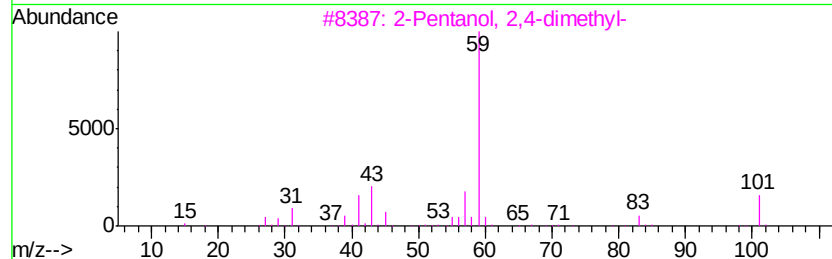
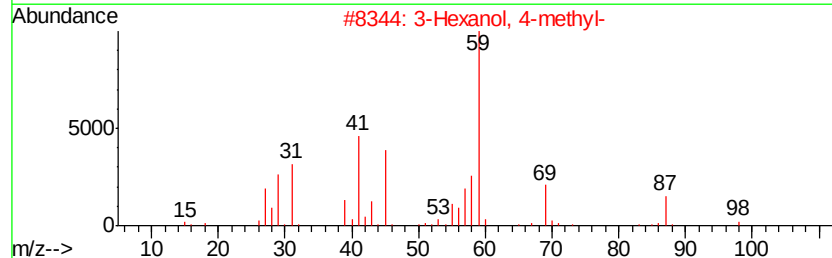
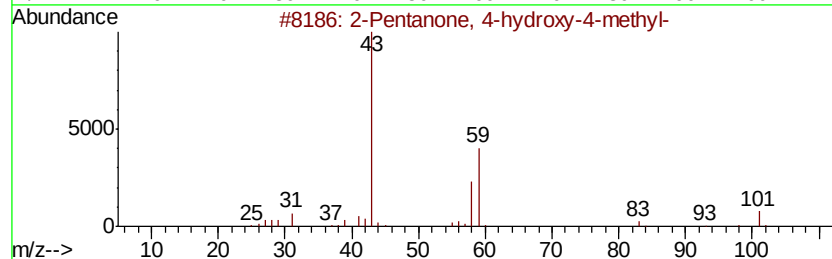
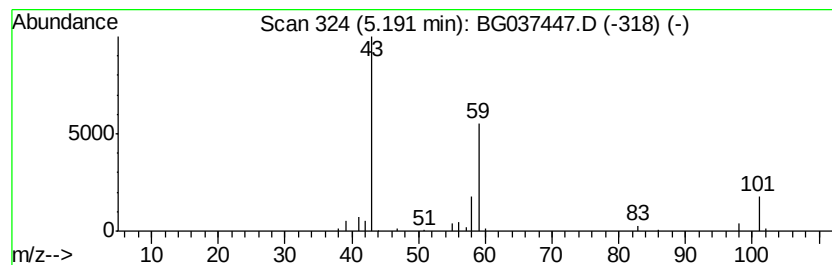
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.19	3.46 ng	63294	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	59	
2	3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	33	
3	2-Pentanol, 2,4-dimethyl-	116	C7H16O	000625-06-9	23	
4	2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	23	
5	Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	17	



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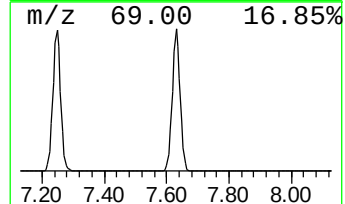
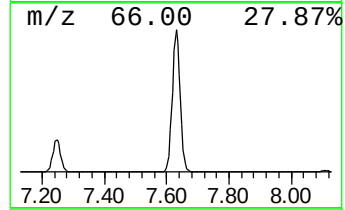
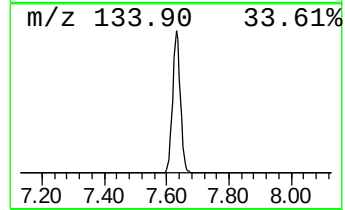
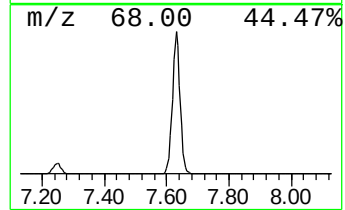
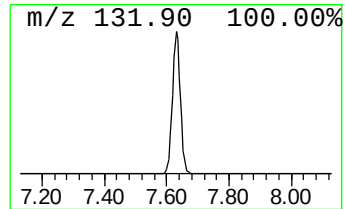
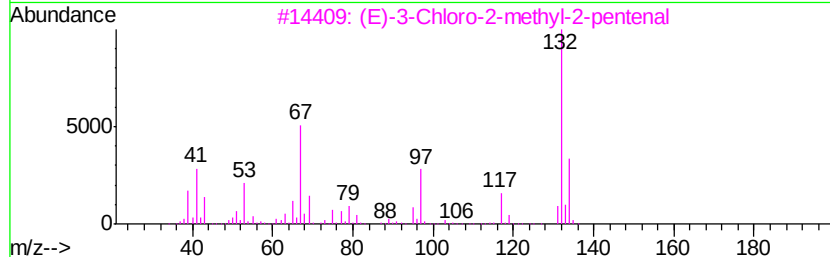
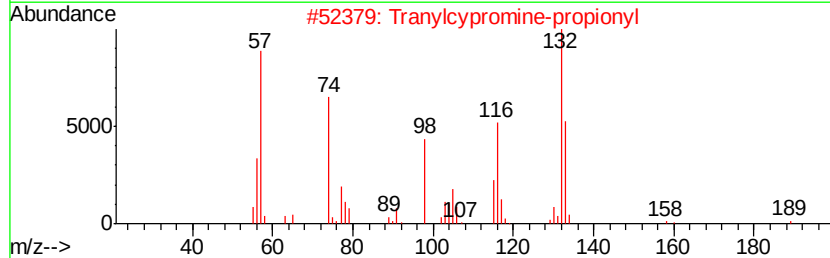
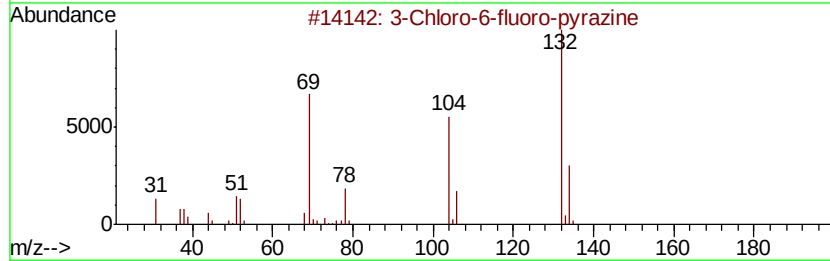
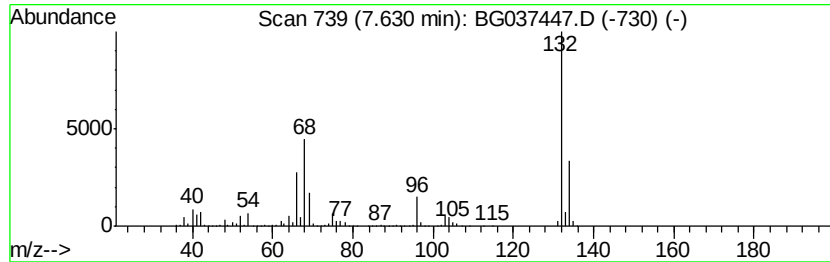
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Peak Number 2 unknown7.63 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.63	90.88 ng	1660500	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		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
3		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	9
4		1,3,5-Trifluorobenzene	132	C6H3F3	000372-38-3	9
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...	5.19	3.5	ng	63294	1	8.11	365444	20.0
unknown7.63	7.63	90.9	ng	1660500	1	8.11	365444	20.0