

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG102518\
 Data File : BG037547.D
 Acq On : 25 Oct 2018 16:20
 Operator : JU/SJ
 Sample : PB114142BL
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB114142BL

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-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	5.180	317	322	330	rVB2	33873	58130	1.72%	0.337%
2	5.644	394	401	412	rBV	782470	1292467	38.28%	7.483%
3	7.248	665	674	687	rBV	829469	1457517	43.17%	8.438%
4	7.630	731	739	748	rBV	757893	1339578	39.68%	7.756%
5	8.106	812	820	828	rBV	160571	285856	8.47%	1.655%
6	8.429	867	875	888	rVB	533134	978298	28.98%	5.664%
7	9.281	1012	1020	1033	rBV	463912	863076	25.56%	4.997%
8	10.926	1292	1300	1309	rVB	236353	438529	12.99%	2.539%
9	13.358	1706	1714	1726	rBV	1375250	2197512	65.09%	12.723%
10	14.739	1941	1949	1963	rBV2	405000	691415	20.48%	4.003%
11	16.220	2193	2201	2221	rBV	962101	1558129	46.15%	9.021%
12	17.483	2408	2416	2427	rBV	593957	900732	26.68%	5.215%
13	19.751	2797	2802	2807	rBV	56381	67425	2.00%	0.390%
14	20.080	2852	2858	2867	rBV	2531968	3376211	100.00%	19.547%
15	20.791	2975	2979	2983	rBV2	39369	48707	1.44%	0.282%
16	21.766	3138	3145	3159	rBV	516699	949576	28.13%	5.498%
17	25.098	3701	3712	3727	rBV	221460	769146	22.78%	4.453%

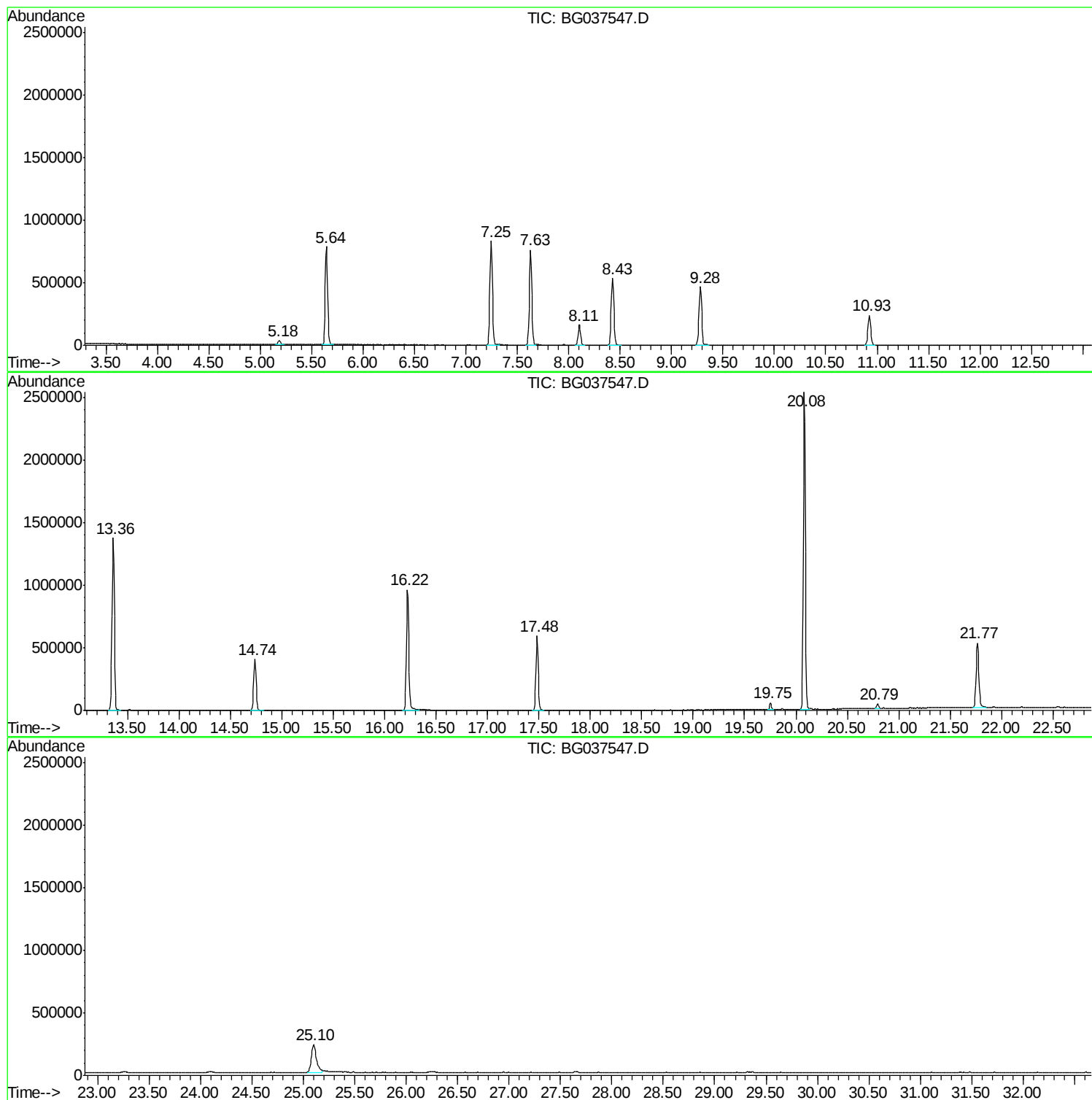
Sum of corrected areas: 17272304

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Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG101818.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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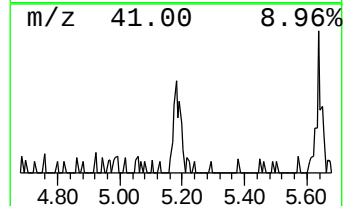
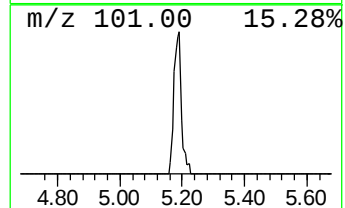
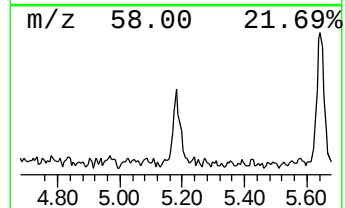
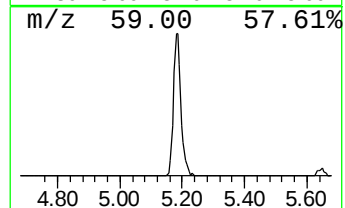
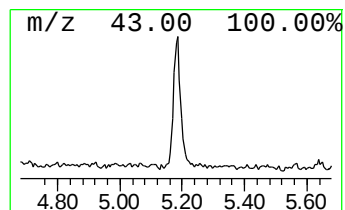
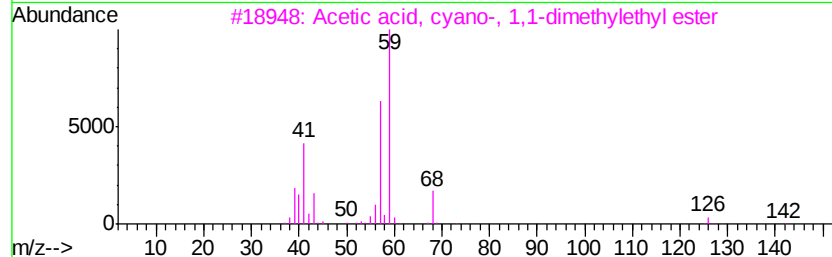
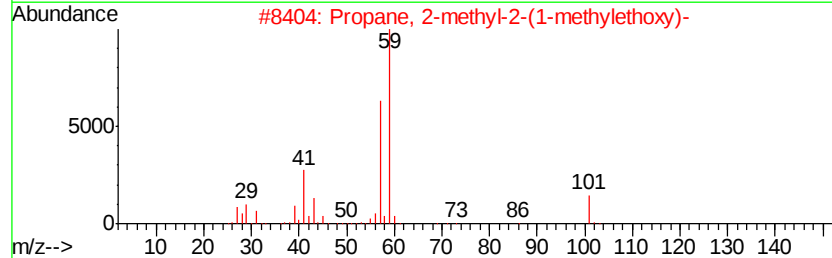
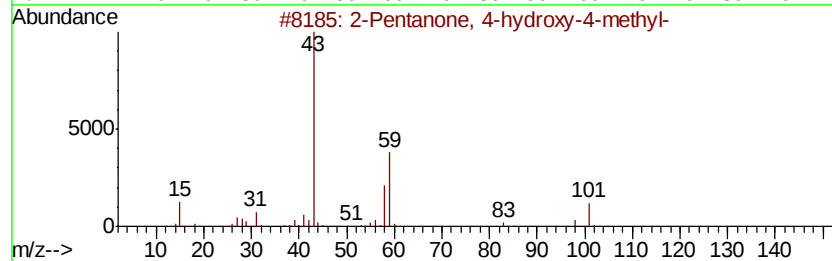
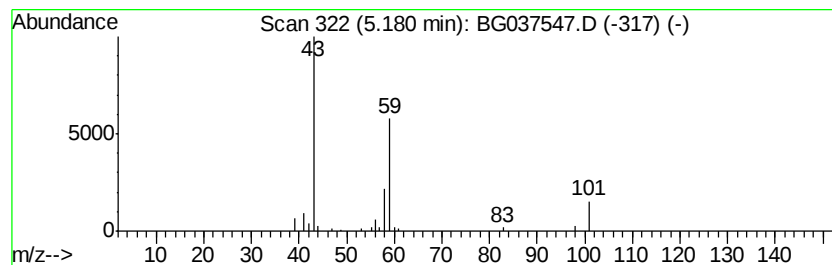
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.18	4.07 ng	58130	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			Propane, 2-methyl-2-(1-methyleth...	116	C7H16O	017348-59-3	28
3			Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	17
4			4-Penten-2-one, 4-methyl-	98	C6H10O	003744-02-3	9
5			Ethyl ether	74	C4H10O	000060-29-7	9



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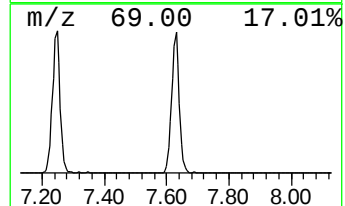
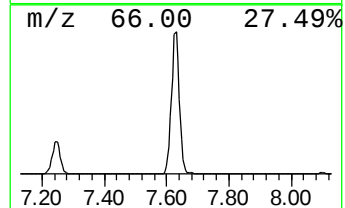
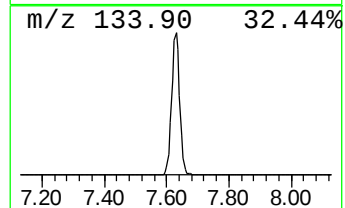
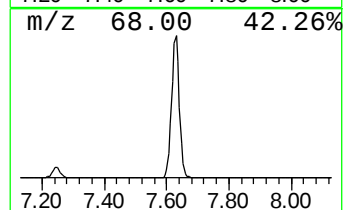
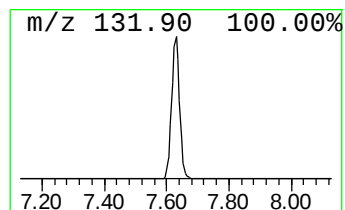
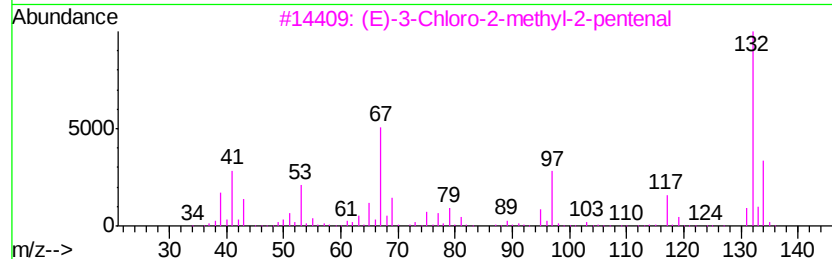
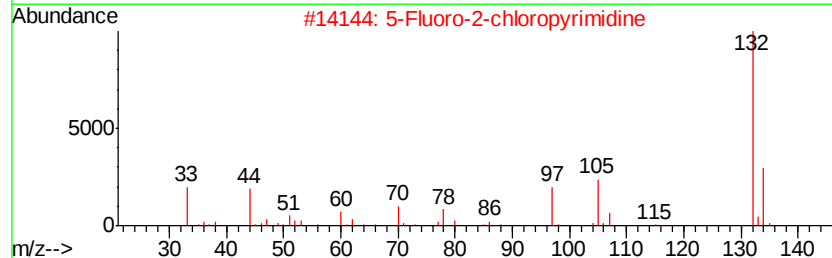
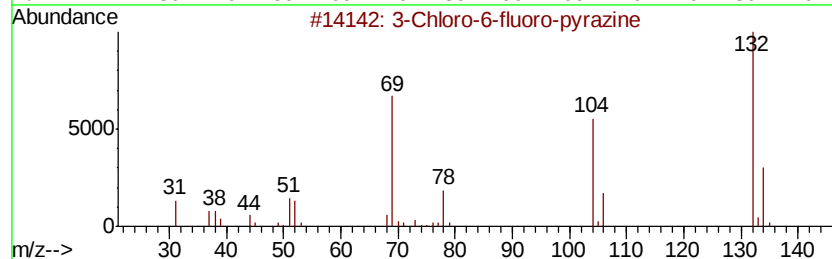
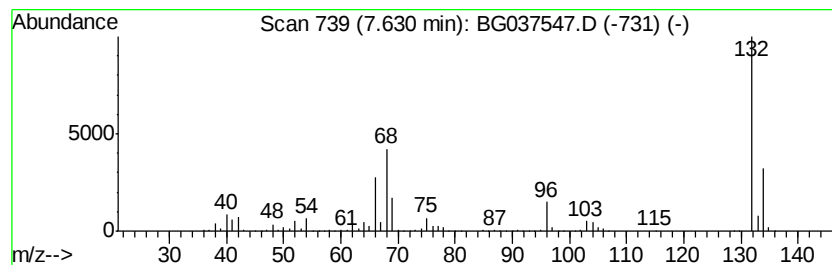
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Peak Number 2 unknown7.63 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.63	93.72 ng	1339580	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		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	16
3		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	12
4		Tranvlcypropine-propionyl	189	C12H15NO	1000123-86-3	10
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...	5.18	4.1	ng	58130	1	8.11	285856	20.0
unknown7.63	7.63	93.7	ng	1339580	1	8.11	285856	20.0