

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG101218\
 Data File : BG037319.D
 Acq On : 13 Oct 2018 10:01
 Operator : JU/SJ
 Sample : J5383-08
 Misc :
 ALS Vial : 27 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 100918-TIPC-F-D-S

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-BG101118.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	5.191	319	324	333	rVB	47548	76271	1.42%	0.360%
2	5.650	395	402	412	rBV	334902	551060	10.22%	2.598%
3	7.254	665	675	685	rBV	207246	408595	7.58%	1.926%
4	7.641	732	741	750	rBV	609990	1086735	20.16%	5.123%
5	8.117	814	822	833	rBV	214480	374483	6.95%	1.765%
6	8.440	868	877	885	rBV	883462	1569189	29.11%	7.397%
7	9.292	1012	1022	1033	rBV	720325	1333250	24.74%	6.285%
8	10.938	1295	1302	1313	rVB	299443	544971	10.11%	2.569%
9	13.376	1709	1717	1727	rBV	2115081	3329326	61.77%	15.695%
10	14.193	1847	1856	1862	rBV	57670	85639	1.59%	0.404%
11	14.751	1942	1951	1960	rBV2	518387	847080	15.72%	3.993%
12	16.231	2197	2203	2218	rBV	1108034	1685360	31.27%	7.945%
13	17.495	2408	2418	2427	rBV	737343	1153782	21.41%	5.439%
14	20.097	2854	2861	2867	rBV	3849623	5389723	100.00%	25.408%
15	21.390	3077	3081	3092	rBV2	55195	95623	1.77%	0.451%
16	21.778	3140	3147	3156	rVB2	715958	1273089	23.62%	6.001%
17	23.505	3436	3441	3449	rVB3	31344	66487	1.23%	0.313%
18	24.151	3544	3551	3557	rBV5	24591	59355	1.10%	0.280%
19	25.127	3706	3717	3732	rVB2	413093	1283039	23.81%	6.048%

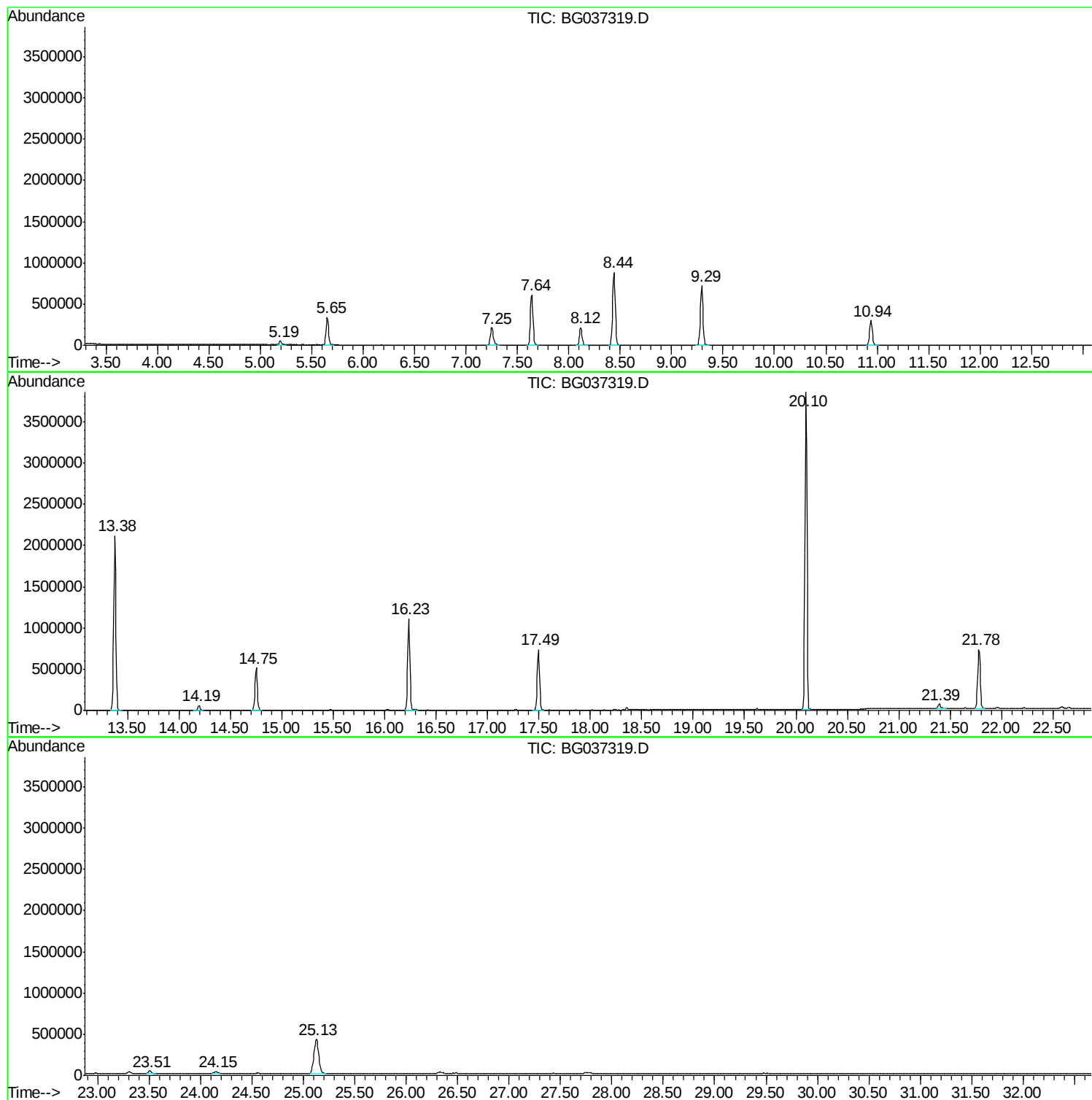
Sum of corrected areas: 21213057

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Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG101118.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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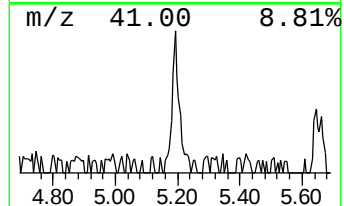
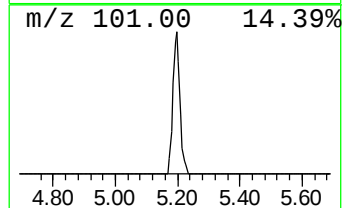
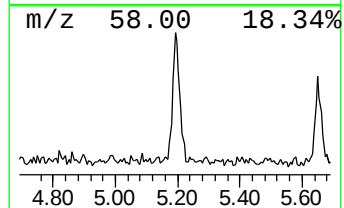
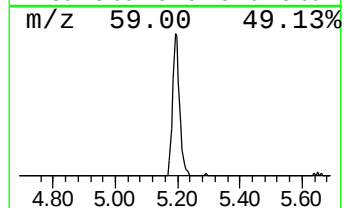
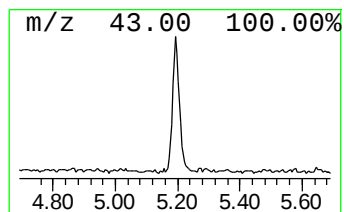
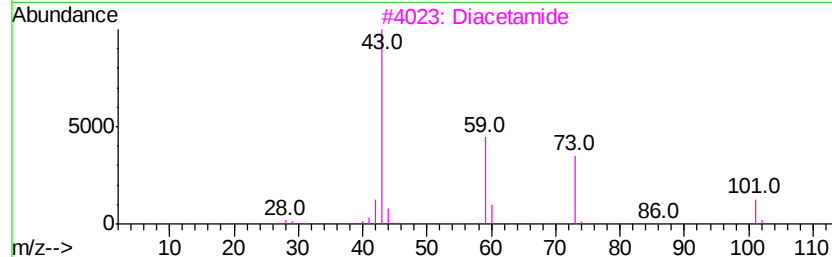
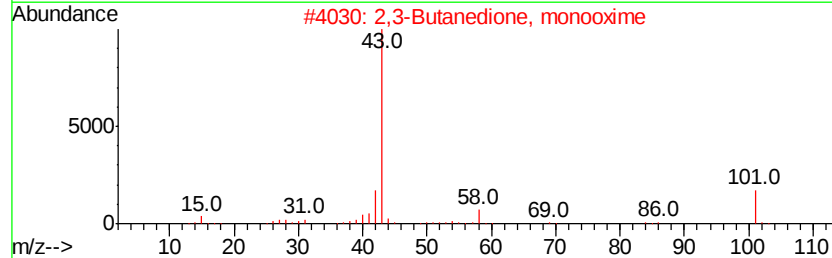
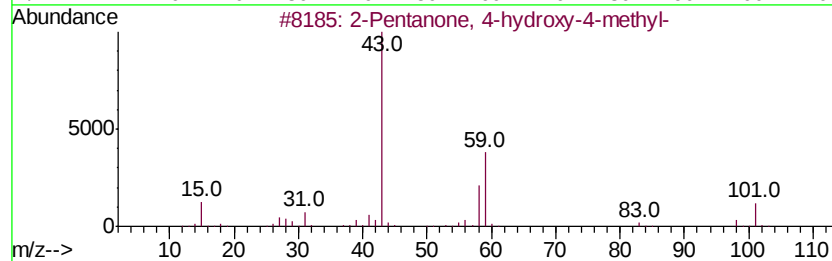
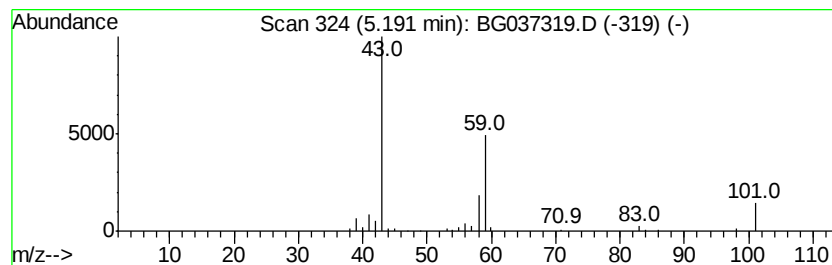
Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG101118.M
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.19	4.07 ng	76271	1,4-Dichlorobenzene-d4	8.12

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50
2			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9
3			Diacetamide	101	C4H7NO2	000625-77-4	9
4			1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	9
5			4-Penten-2-one, 4-methyl-	98	C6H10O	003744-02-3	7



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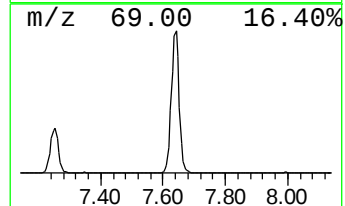
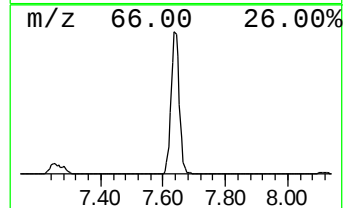
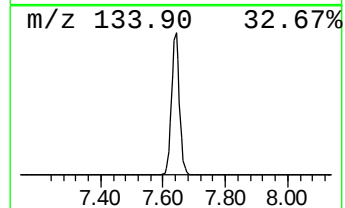
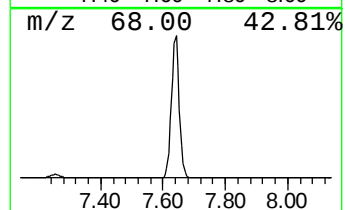
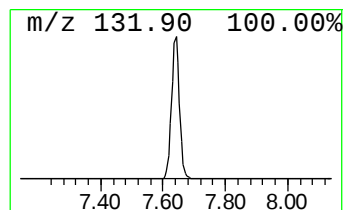
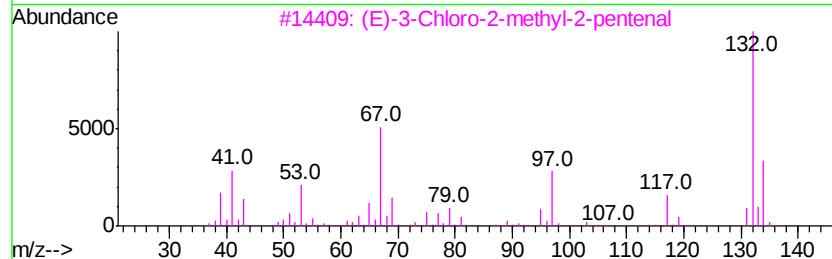
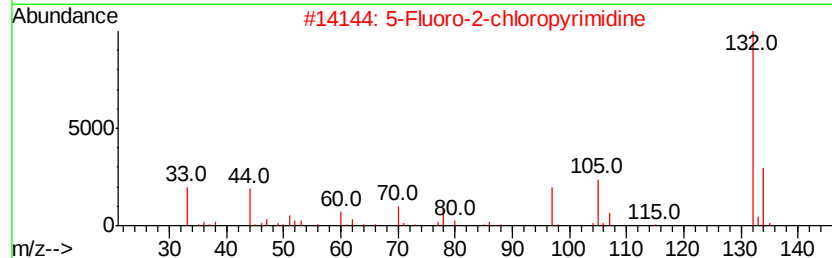
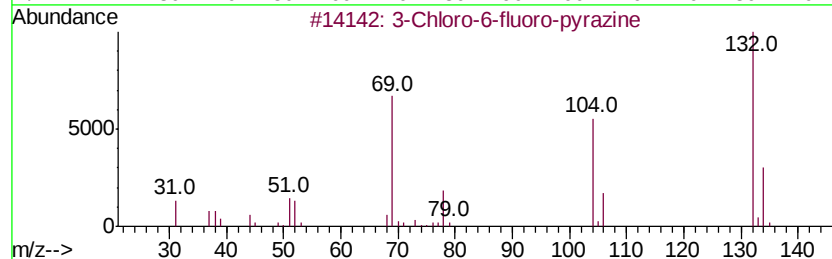
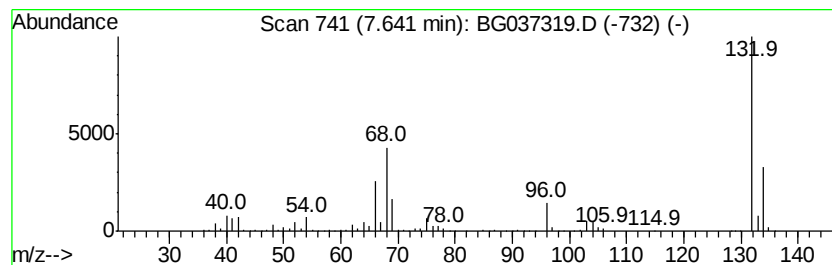
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Peak Number 2 unknown7.64 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.64	58.04 ng	1086740	1,4-Dichlorobenzene-d4	8.12

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	16
4		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
5		3H-1,2-Dithiol-3-one, 5-methyl-	132	C4H4OS2	003620-08-4	9



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TIC Top Hit name	RT	EstConc	Units	Response	---Internal Standard---			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.19	4.1	ng	76271	1	8.12	374483	20.0
unknown7.64	7.64	58.0	ng	1086740	1	8.12	374483	20.0