

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG101218\
 Data File : BG037312.D
 Acq On : 13 Oct 2018 5:27
 Operator : JU/SJ
 Sample : J5383-01
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 100918-SIDI-F-D-B

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	5.196	319	325	333	rBV	67630	112799	2.06%	0.532%
2	5.649	396	402	412	rVB	352843	594676	10.88%	2.802%
3	7.253	666	675	686	rBV	239496	443133	8.11%	2.088%
4	7.640	733	741	752	rBV	643316	1137863	20.83%	5.362%
5	8.116	815	822	835	rVB	203439	357838	6.55%	1.686%
6	8.440	868	877	886	rVB	872882	1578049	28.88%	7.436%
7	9.291	1014	1022	1034	rBV	726338	1351984	24.75%	6.371%
8	10.937	1295	1302	1311	rBV	269756	518487	9.49%	2.443%
9	13.375	1709	1717	1725	rBV	2119896	3379894	61.86%	15.926%
10	14.750	1943	1951	1959	rVB2	489268	817433	14.96%	3.852%
11	16.236	2196	2204	2217	rBV2	1091258	1721399	31.51%	8.111%
12	17.494	2411	2418	2432	rBV	679028	1102019	20.17%	5.193%
13	18.357	2560	2565	2574	rBV2	54921	92081	1.69%	0.434%
14	20.096	2855	2861	2875	rBV	3855041	5463617	100.00%	25.744%
15	21.783	3141	3148	3157	rVB	712369	1190808	21.80%	5.611%
16	23.510	3436	3442	3451	rBV2	49750	102535	1.88%	0.483%
17	24.151	3546	3551	3561	rVB3	27853	64773	1.19%	0.305%
18	25.126	3706	3717	3731	rBV2	386056	1193109	21.84%	5.622%

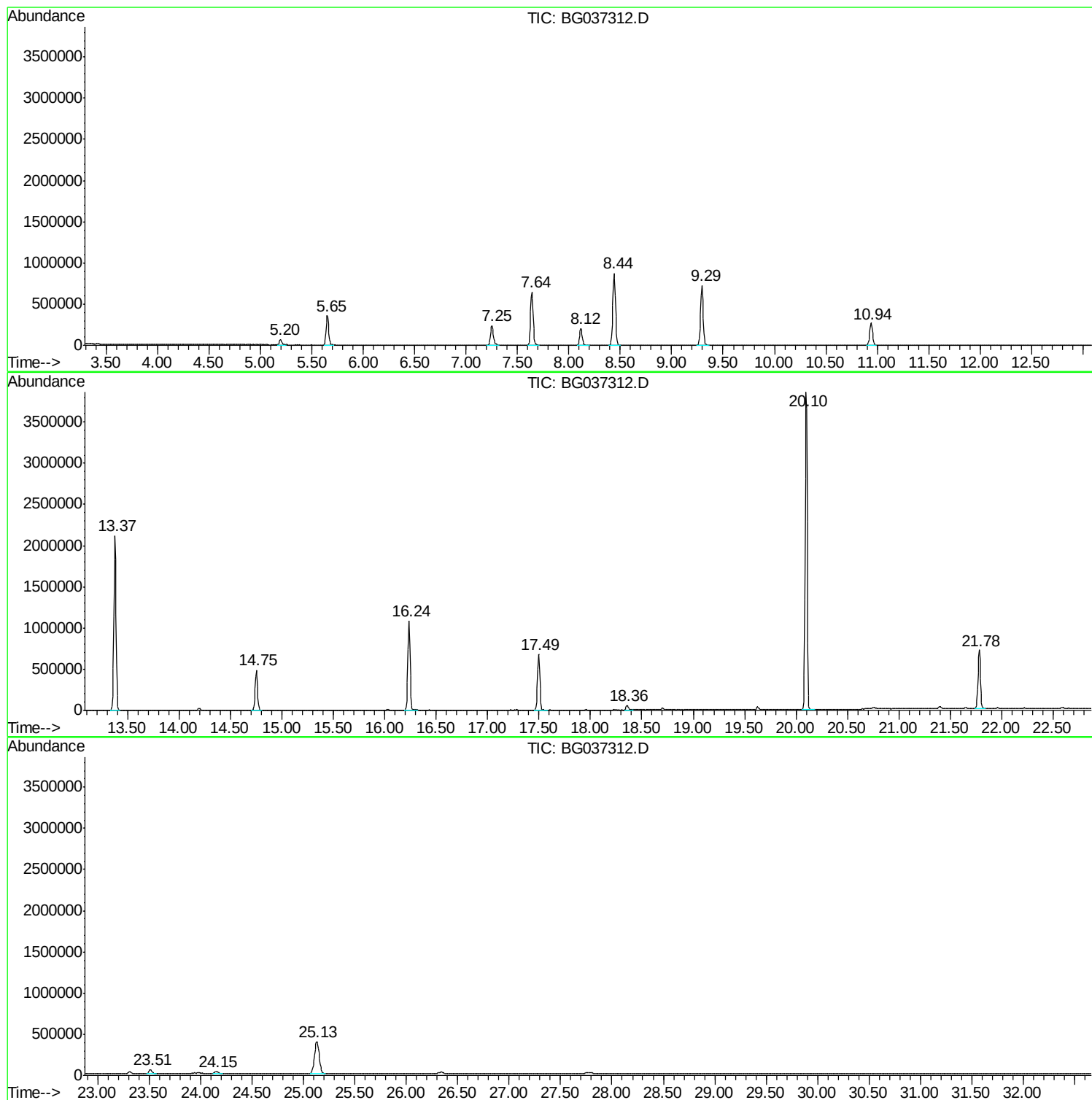
Sum of corrected areas: 21222497

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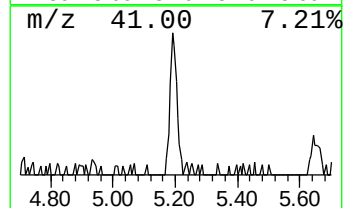
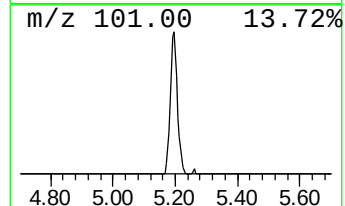
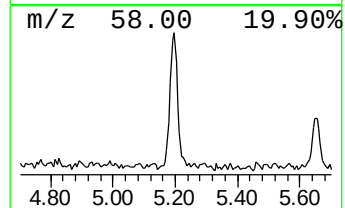
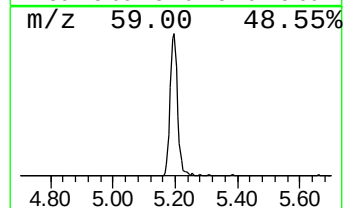
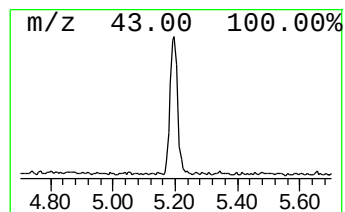
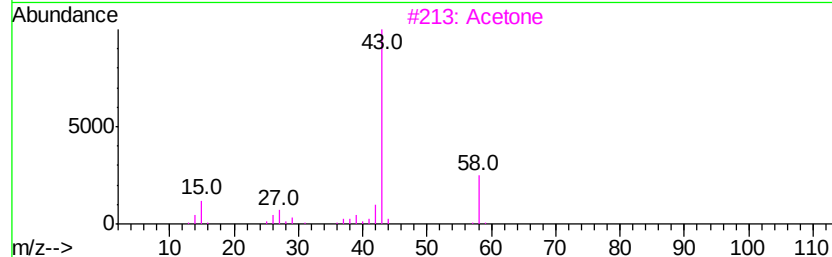
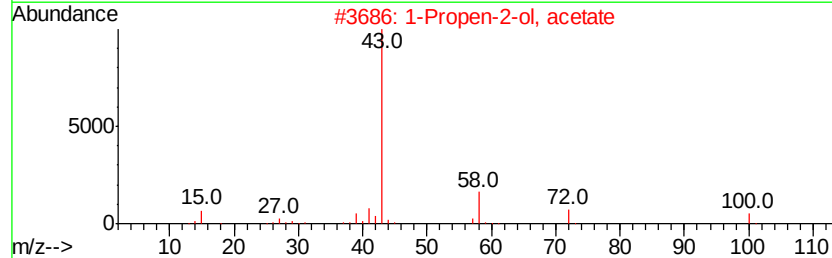
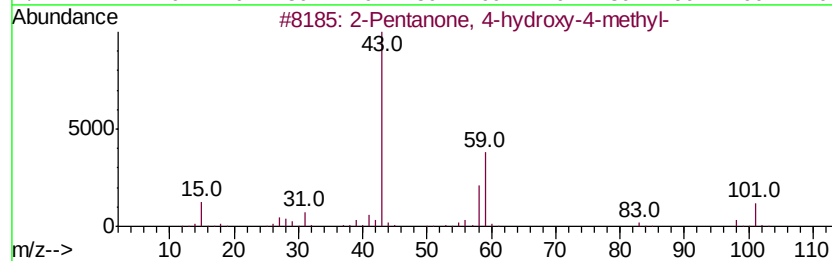
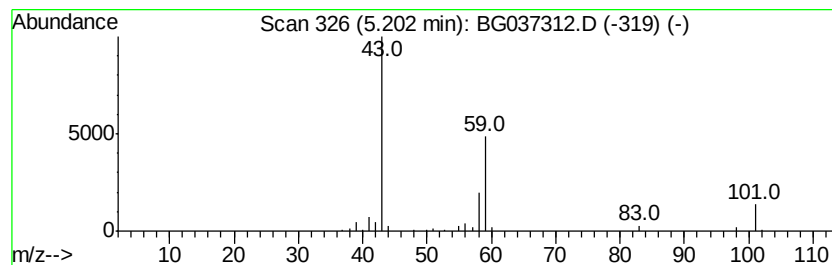
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.20	6.30 ng	112799	1,4-Dichlorobenzene-d4	8.12

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	78
2		1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	12
3		Acetone	58	C3H6O	000067-64-1	9
4		Diacetamide	101	C4H7NO2	000625-77-4	9
5		2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9



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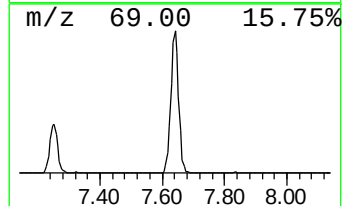
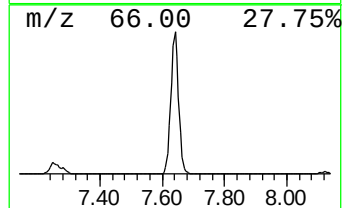
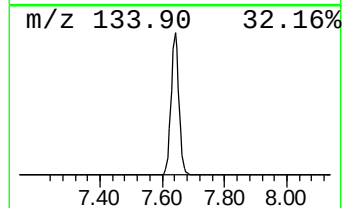
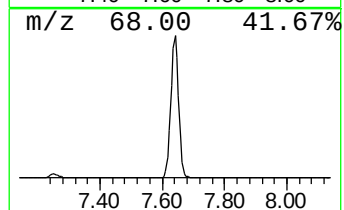
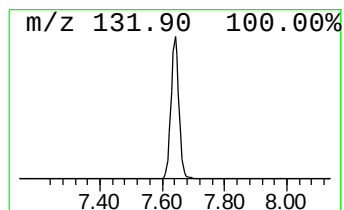
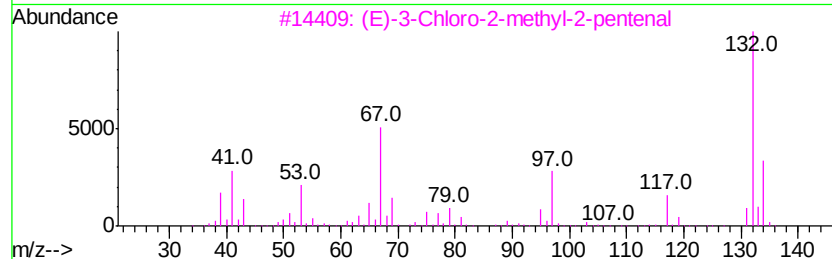
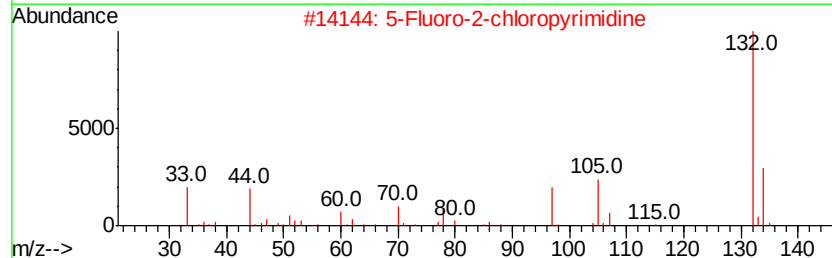
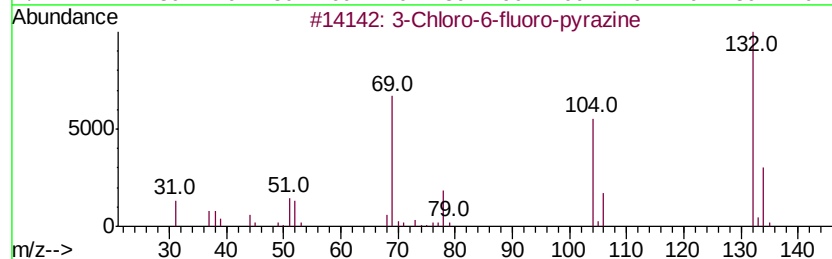
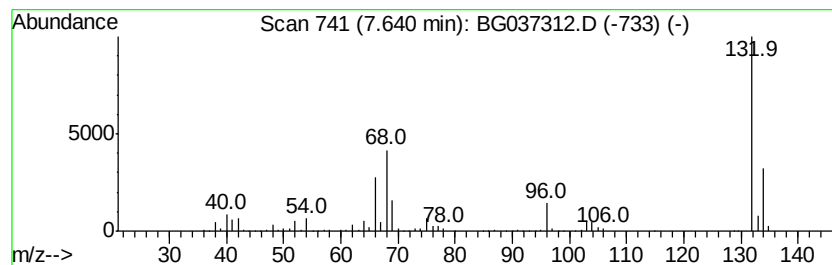
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Peak Number 2 unknown7.64 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.64	63.60 ng	1137860	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	12
4		Tranvlcypropine-propionyl	189	C12H15NO	1000123-86-3	10
5		2H-Cyclopenta[d]pyridazine, 2-me...	132	C8H8N2	022291-85-6	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.20	6.3	ng	112799	1	8.12	357838	20.0
unknown7.64	7.64	63.6	ng	1137860	1	8.12	357838	20.0