

Data Path : Z:\HPCHEM1\BNA G\DATA\BG112217\
 Data File : BG031193.D
 Acq On : 22 Nov 2017 17:25
 Operator : SJ/JU
 Sample : I6525-03
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 T71-MW7

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:\HPCHEM1\BNA_G\METHODS\8270-BG111017.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	5.212	321	327	336	rBV	51765	78903	1.42%	0.372%
2	5.694	403	409	430	rBV	398897	696332	12.50%	3.281%
3	7.304	676	683	695	rBV	231497	448662	8.05%	2.114%
4	7.668	737	745	757	rBV	778083	1412399	25.35%	6.656%
5	8.138	818	825	834	rBV	172384	304213	5.46%	1.434%
6	8.461	872	880	892	rBV	775757	1400071	25.12%	6.598%
7	9.302	1015	1023	1041	rBV	551576	1067780	19.16%	5.032%
8	10.953	1295	1304	1322	rBV	215315	458476	8.23%	2.161%
9	13.379	1709	1717	1741	rBV	1962276	3124933	56.08%	14.726%
10	14.754	1945	1951	1965	rBV2	393766	704860	12.65%	3.322%
11	16.240	2197	2204	2223	rBV	1455660	2361253	42.37%	11.127%
12	17.498	2411	2418	2428	rBV	603161	993964	17.84%	4.684%
13	20.089	2852	2859	2875	rBV	4164013	5572660	100.00%	26.260%
14	21.769	3137	3145	3157	rBV	752653	1342543	24.09%	6.327%
15	25.065	3695	3706	3722	rBV2	408853	1253685	22.50%	5.908%

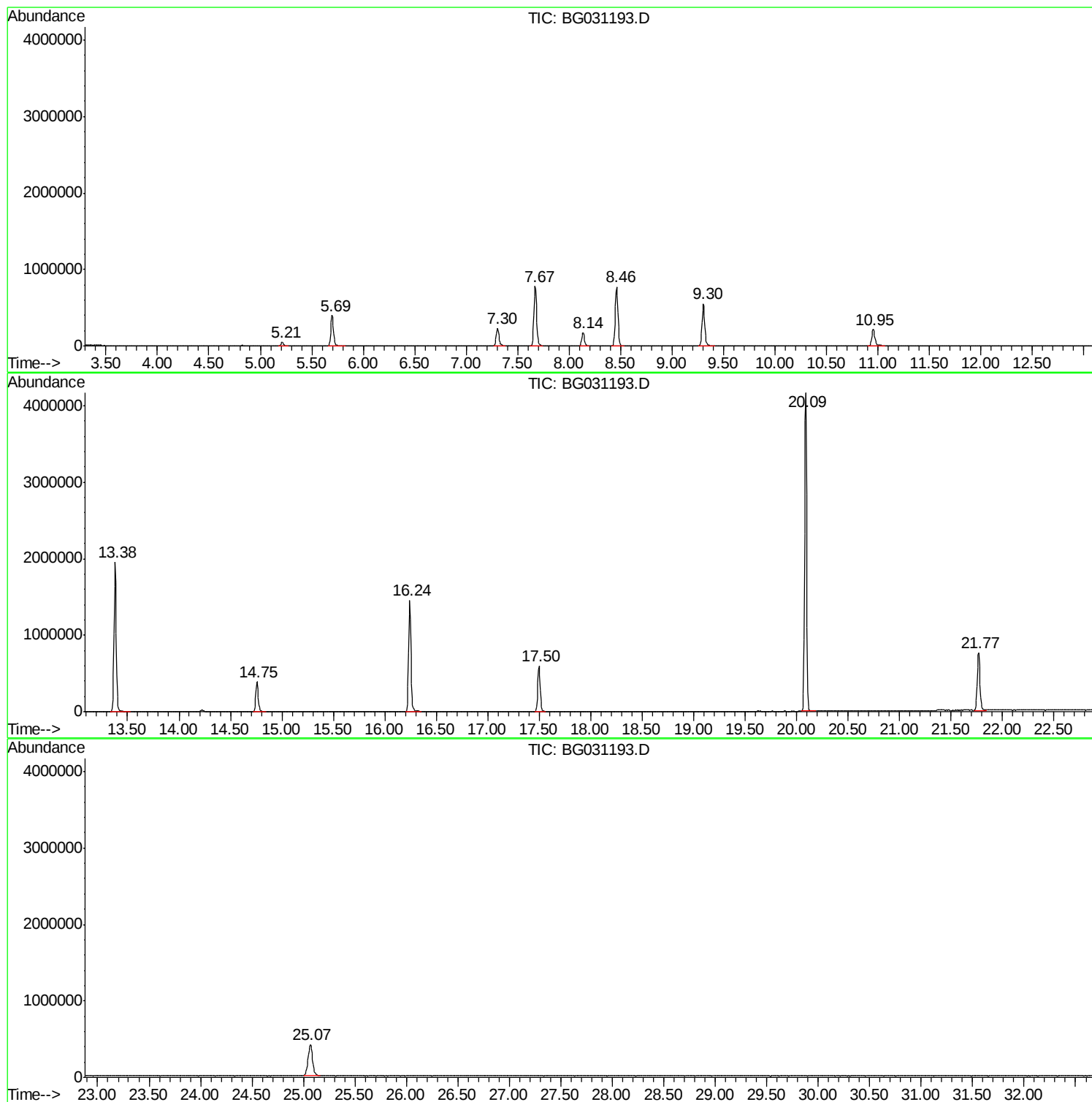
Sum of corrected areas: 21220734

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Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG111017.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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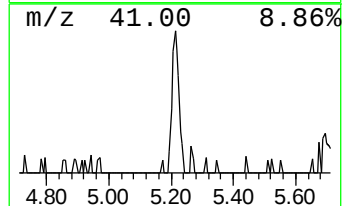
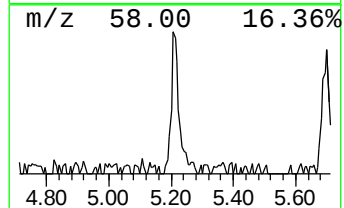
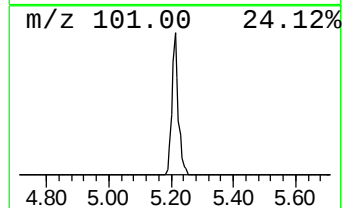
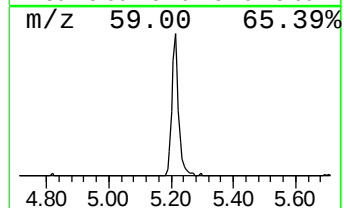
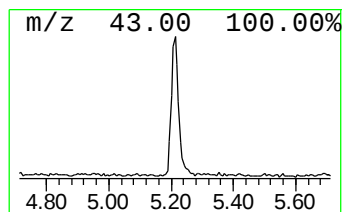
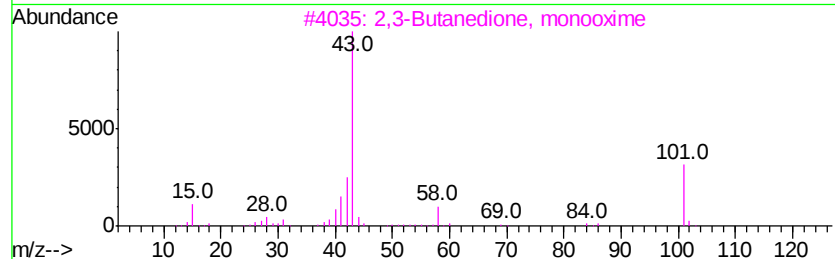
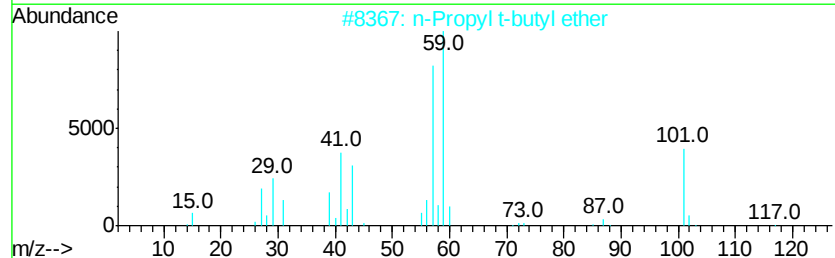
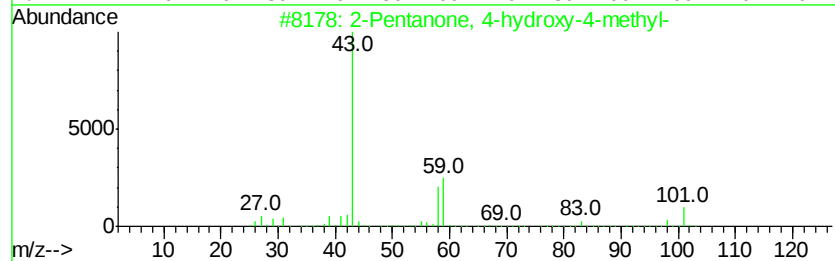
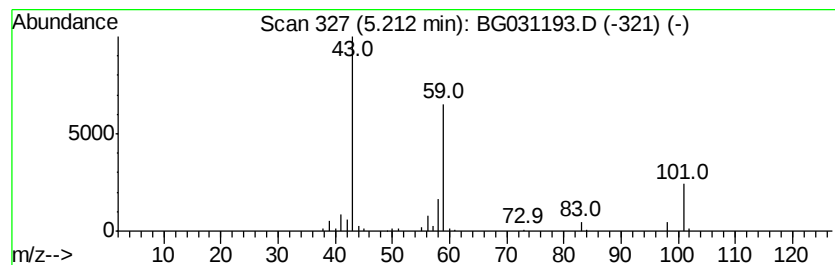
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.21	5.19 ng	78903	1,4-Dichlorobenzene-d4	8.13

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	64
2			n-Propyl t-butyl ether	116	C7H16O	1000334-81-9	33
3			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	16
4			Acetone	58	C3H6O	000067-64-1	14
5			5-Hexen-2-one	98	C6H10O	000109-49-9	9



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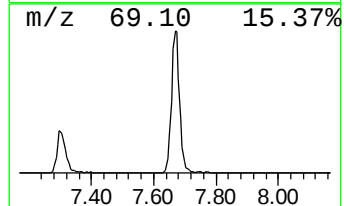
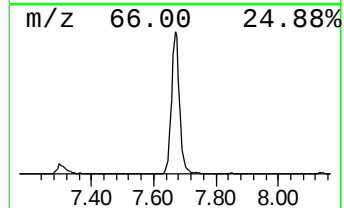
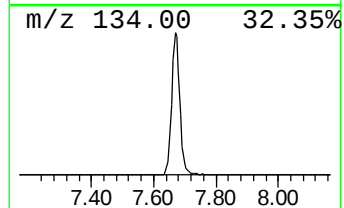
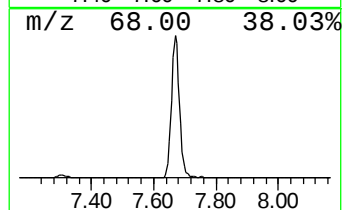
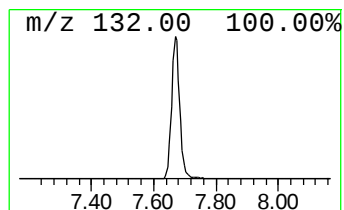
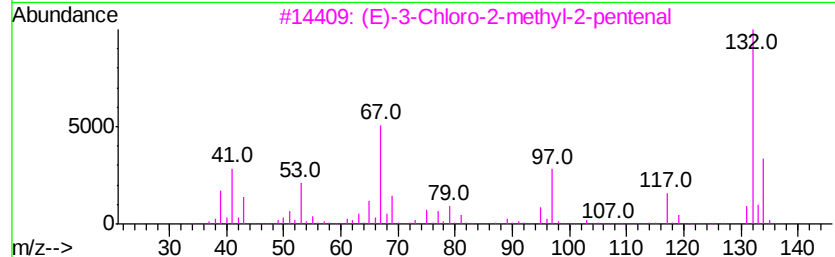
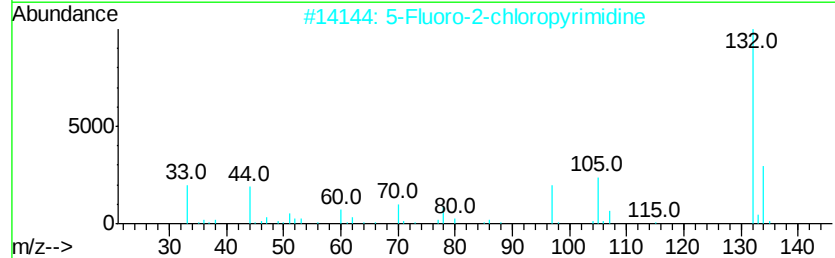
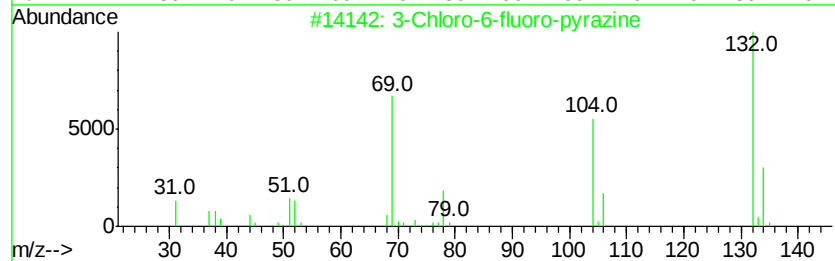
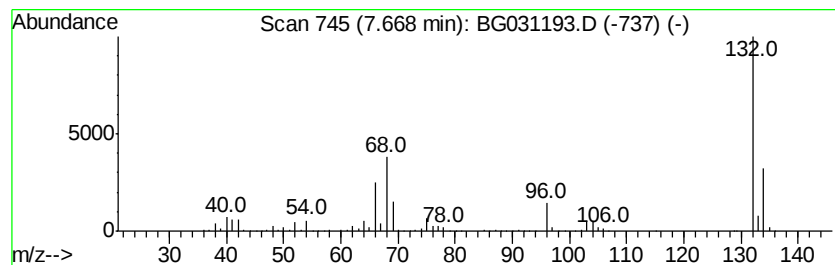
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Peak Number 2 unknown7.67 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.67	92.86 ng	1412400	1,4-Dichlorobenzene-d4	8.13

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	35
2		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	25
3		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	17
4		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	12
5		5-Aminoindole	132	C8H8N2	005192-03-0	12



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.21	5.2	ng	78903	1	8.13	304213	20.0
unknown7.67	7.67	92.9	ng	1412400	1	8.13	304213	20.0