

Data Path : Z:\HPCHEM1\BNA G\DATA\BG120617\
 Data File : BG031372.D
 Acq On : 6 Dec 2017 15:50
 Operator : SJ/JU
 Sample : PB104763BL
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB104763BL

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	5.204	321	326	334	rBV	61213	96909	1.67%	0.399%
2	5.686	402	408	426	rBV	967603	1689890	29.19%	6.960%
3	7.296	674	682	704	rBV	974162	1776599	30.69%	7.318%
4	7.660	736	744	757	rBV	1018229	1792844	30.97%	7.385%
5	8.125	816	823	836	rBV	189345	340099	5.87%	1.401%
6	8.454	870	879	890	rBV	778966	1422452	24.57%	5.859%
7	9.300	1015	1023	1044	rBV	527607	1090578	18.84%	4.492%
8	10.945	1296	1303	1313	rBV	215337	468137	8.09%	1.928%
9	13.371	1709	1716	1727	rBV	1778335	2990458	51.65%	12.317%
10	14.752	1944	1951	1966	rBV2	402008	746182	12.89%	3.073%
11	16.239	2197	2204	2229	rBV	1214989	2219484	38.34%	9.142%
12	17.496	2411	2418	2434	rBV2	613354	1074261	18.55%	4.425%
13	20.081	2852	2858	2881	rBV	4213999	5789693	100.00%	23.847%
14	21.767	3138	3145	3164	rBV	789830	1433634	24.76%	5.905%
15	25.064	3694	3706	3726	rBV2	411031	1347255	23.27%	5.549%

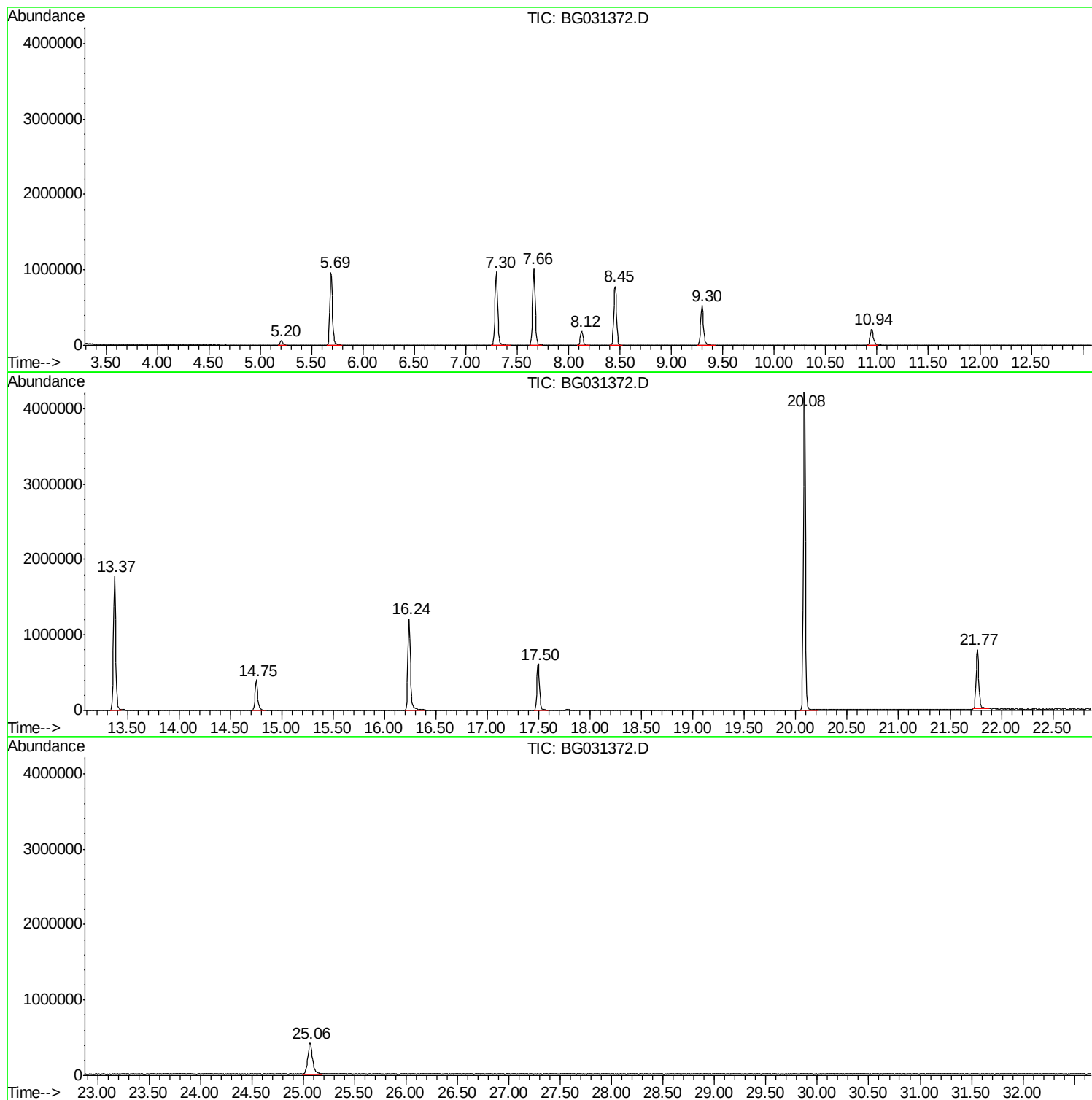
Sum of corrected areas: 24278475

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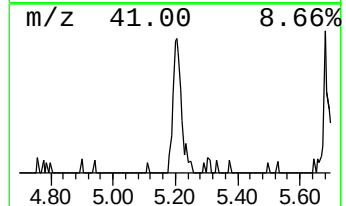
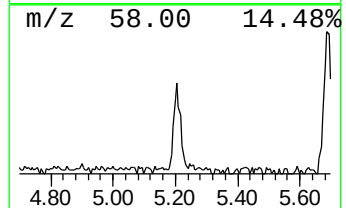
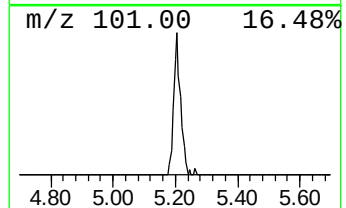
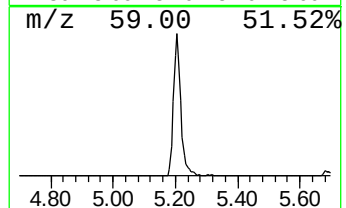
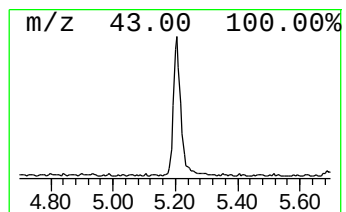
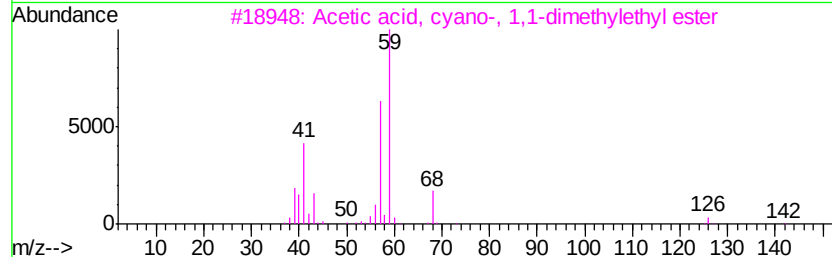
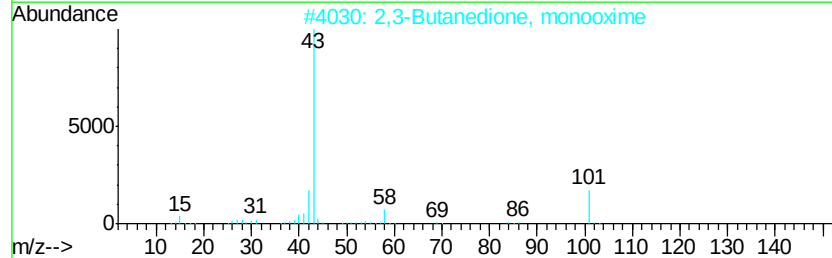
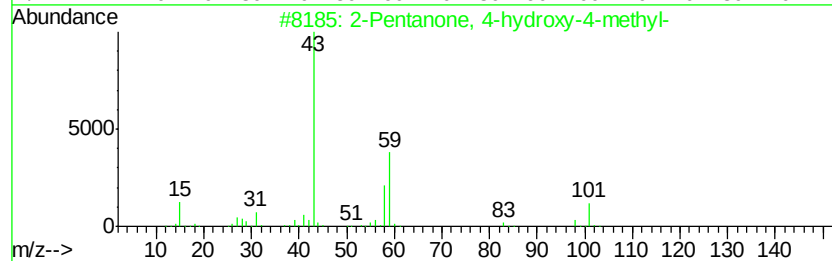
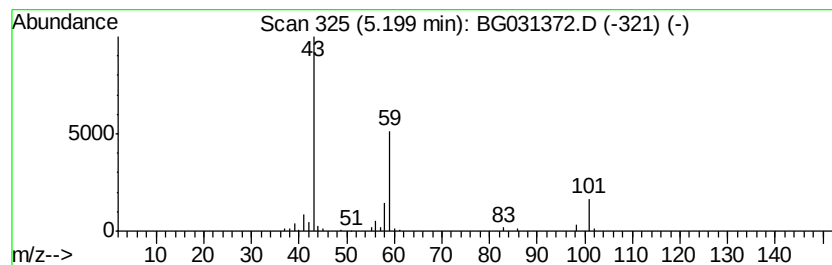
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.20	5.70 ng	96909	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	56
2	2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	38
3	Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	23
4	1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	10
5	Butane	58	C4H10	000106-97-8	9



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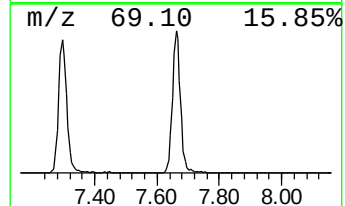
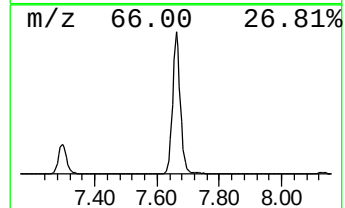
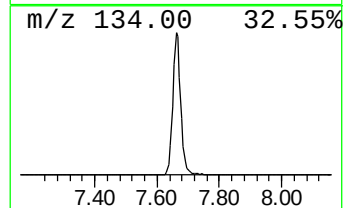
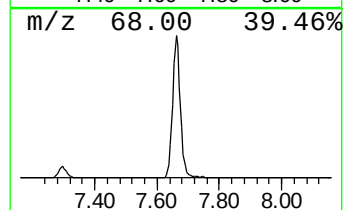
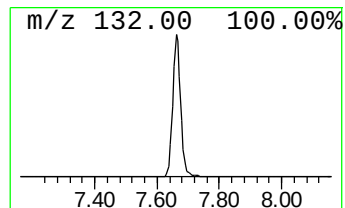
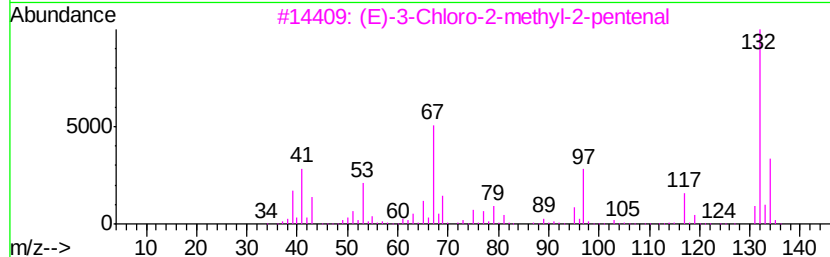
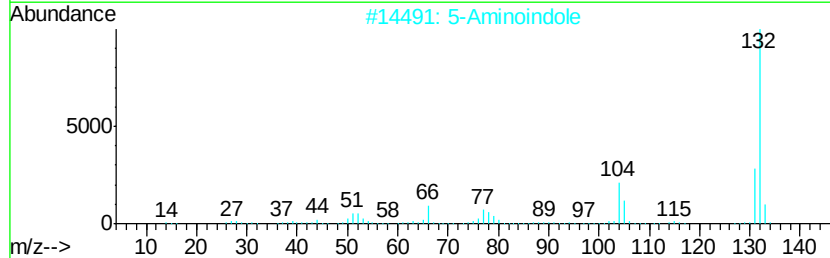
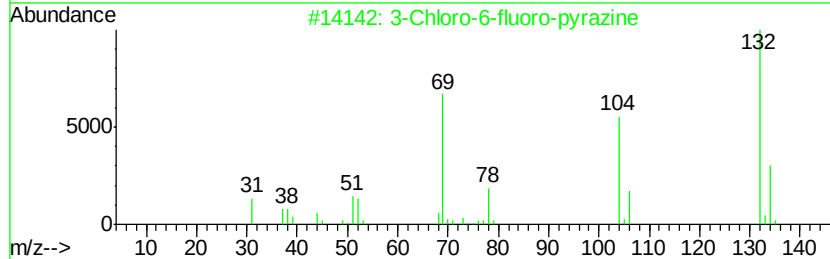
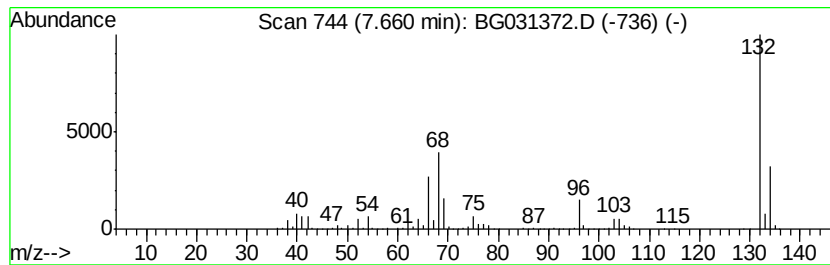
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Peak Number 2 unknown7.66 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.66	105.43 ng	1792840	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	27
2		5-Aminoindole	132	C8H8N2	005192-03-0	22
3		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	16
4		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	12
5		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.20	5.7	ng	96909	1	8.13	340099	20.0
unknown7.66	7.66	105.4	ng	1792840	1	8.13	340099	20.0