

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA_G\DATA\BG011818\
 Data File : BG032141.D
 Acq On : 18 Jan 2018 18:16
 Operator : SJ/JU
 Sample : PB105758BL
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB105758BL

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-BG011218.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	5.662	398	404	411	rBV	29780	50950	1.17%	0.299%
2	6.168	483	490	499	rBV	591406	1032265	23.62%	6.064%
3	7.848	767	776	787	rBV	588263	1062786	24.32%	6.243%
4	8.253	837	845	856	rBV	641094	1143860	26.18%	6.719%
5	8.741	921	928	939	rBV	130371	231680	5.30%	1.361%
6	9.076	977	985	997	rBV	515293	956322	21.88%	5.618%
7	9.940	1123	1132	1145	rBV	342309	654110	14.97%	3.842%
8	11.626	1411	1419	1431	rVB	169608	319533	7.31%	1.877%
9	14.005	1816	1824	1837	rBV	1297405	2061667	47.18%	12.111%
10	15.380	2050	2058	2070	rVB2	291712	486470	11.13%	2.858%
11	16.855	2301	2309	2322	rBV	1124233	1794636	41.07%	10.542%
12	18.112	2516	2523	2532	rBV2	460690	715909	16.38%	4.205%
13	20.651	2949	2955	2964	rBV	3241958	4369941	100.00%	25.670%
14	22.448	3252	3261	3272	rVB2	586957	1080267	24.72%	6.346%
15	26.150	3878	3891	3904	rBV2	324249	1063163	24.33%	6.245%

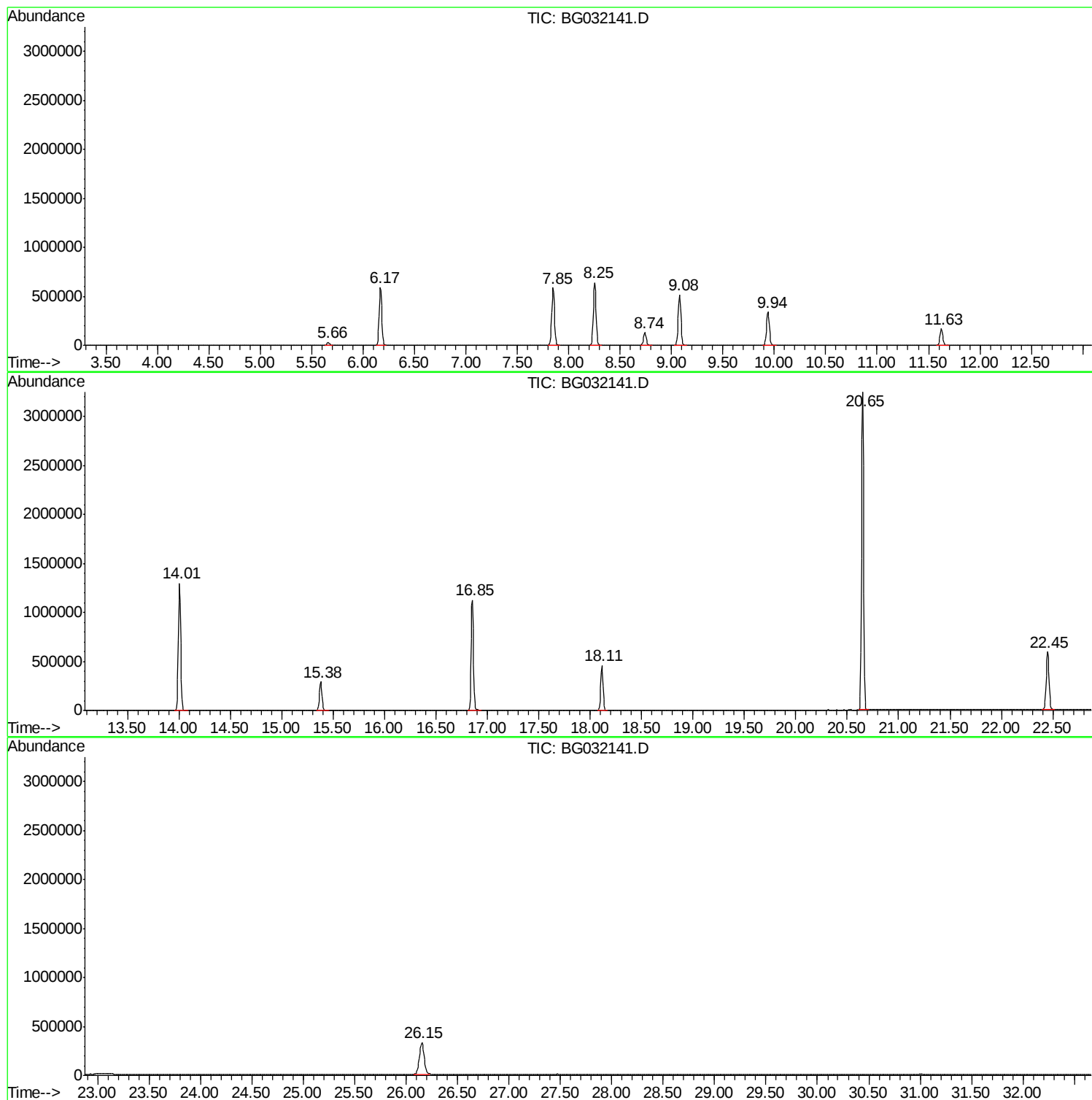
Sum of corrected areas: 17023559

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Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG011218.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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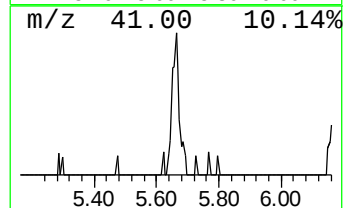
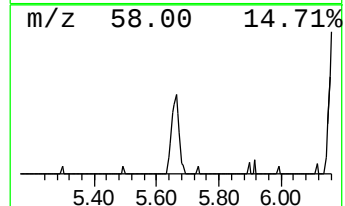
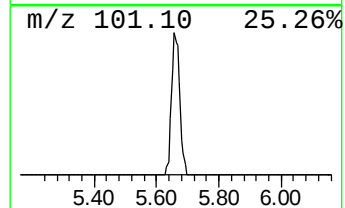
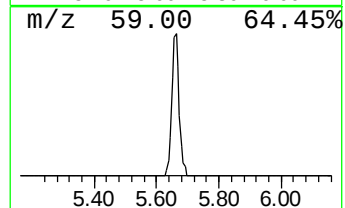
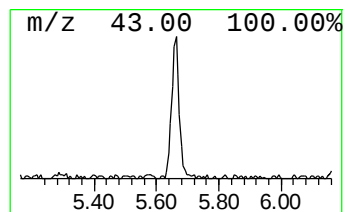
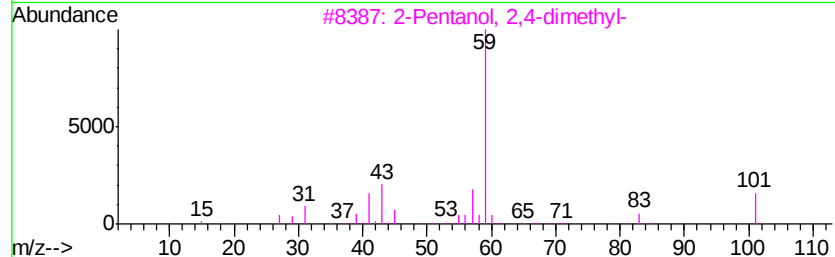
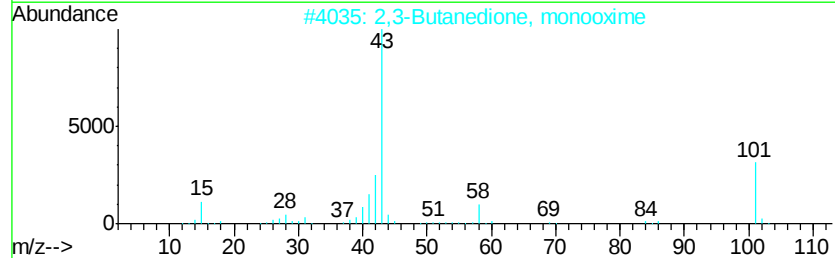
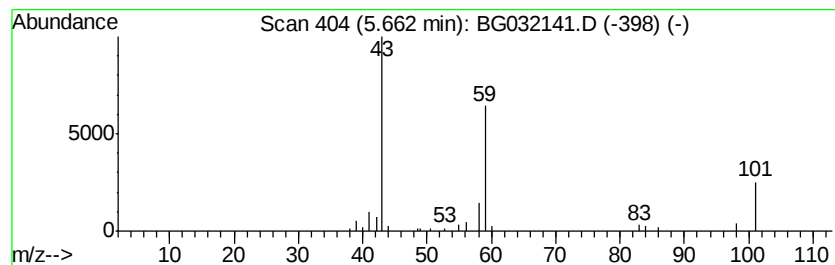
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.66	4.40 ng	50950	1,4-Dichlorobenzene-d4	8.74

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	45
2			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	38
3			2-Pentanol, 2,4-dimethyl-	116	C7H16O	000625-06-9	17
4			1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	9
5			2-Propanone, 1-(1-methylethoxy)-	116	C6H12O2	042781-12-4	9



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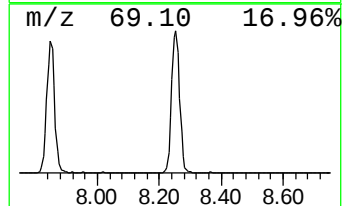
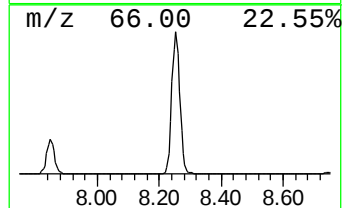
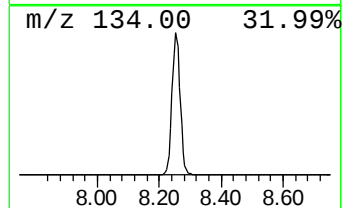
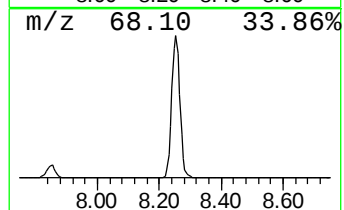
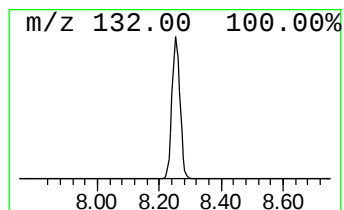
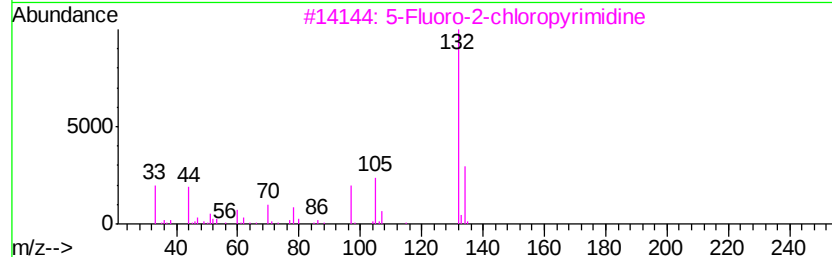
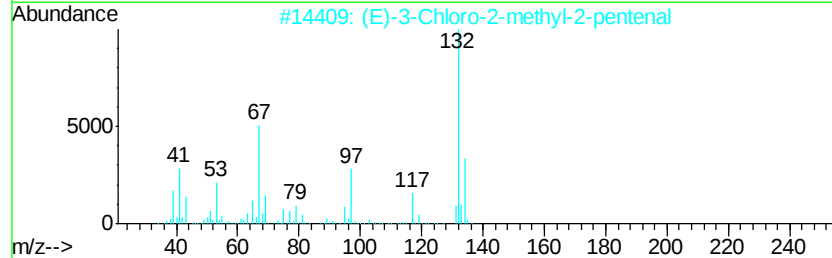
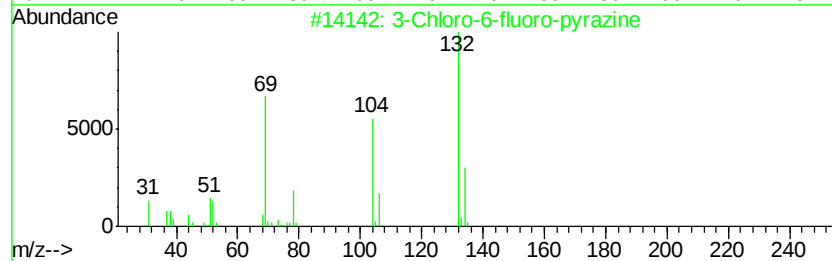
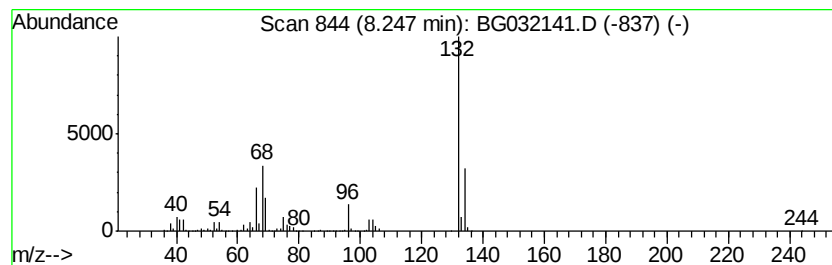
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Peak Number 2 unknown8.25 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.25	98.74 ng	1143860	1,4-Dichlorobenzene-d4	8.74

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	35
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	35
3		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	25
4		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	12
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...	5.66	4.4	ng	50950	1	8.74	231680	20.0
unknown8.25	8.25	98.7	ng	1143860	1	8.74	231680	20.0