

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA G\DATA\BG091017\
 Data File : BG028617.D
 Acq On : 9 Sep 2017 16:21
 Operator : SJ/JU
 Sample : PB102150BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB102150BL

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-BG090717.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	5.441	310	315	322	rBV	50610	79632	2.23%	0.547%
2	5.906	387	394	410	rBV	631171	1075808	30.18%	7.395%
3	7.492	656	664	678	rBV	603141	1095481	30.73%	7.530%
4	7.868	721	728	740	rBV	598747	1063686	29.84%	7.311%
5	8.338	801	808	815	rBV2	106524	192174	5.39%	1.321%
6	8.661	854	863	875	rBV	440698	806036	22.61%	5.540%
7	9.501	998	1006	1018	rBV	342589	662515	18.59%	4.554%
8	11.158	1279	1288	1296	rBV	135570	264503	7.42%	1.818%
9	13.567	1691	1698	1709	rBV	1023375	1634107	45.84%	11.232%
10	14.948	1926	1933	1941	rBV2	228390	388332	10.89%	2.669%
11	16.434	2179	2186	2202	rBV	981862	1599824	44.88%	10.997%
12	17.698	2394	2401	2408	rBV2	356188	589965	16.55%	4.055%
13	20.283	2834	2841	2851	rBV	2669194	3564546	100.00%	24.502%
14	22.028	3130	3138	3148	rBV	435739	780275	21.89%	5.363%
15	25.530	3720	3734	3747	rVB	227299	751262	21.08%	5.164%

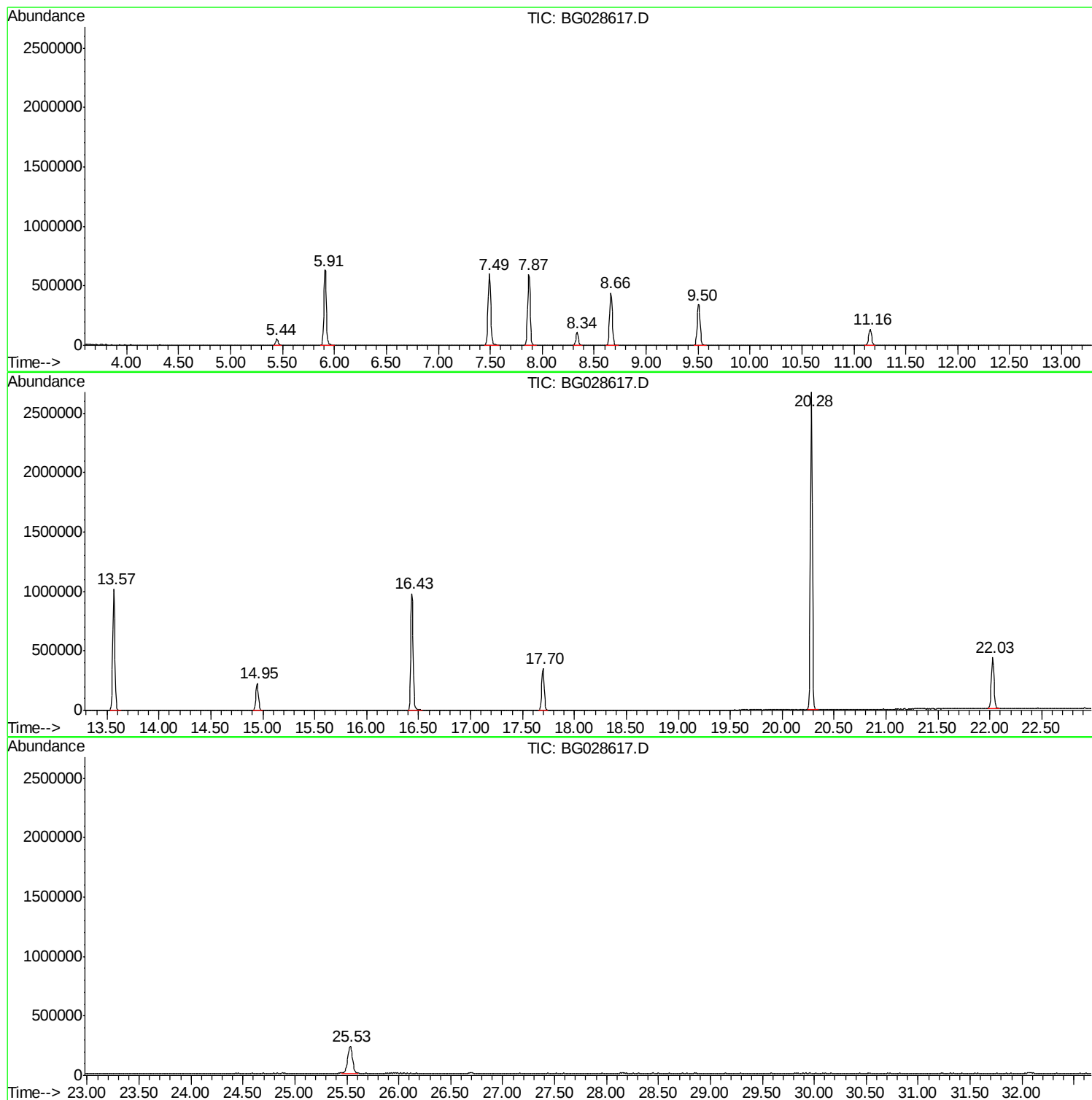
Sum of corrected areas: 14548146

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Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG090717.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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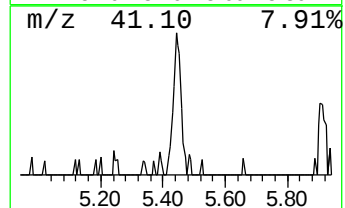
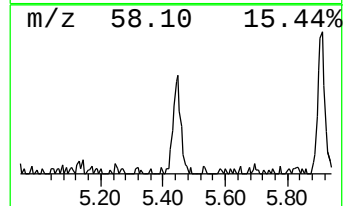
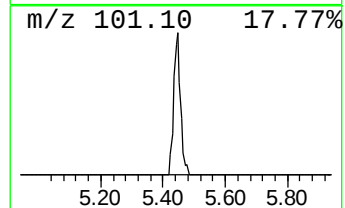
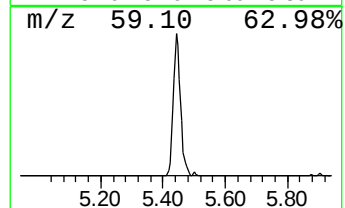
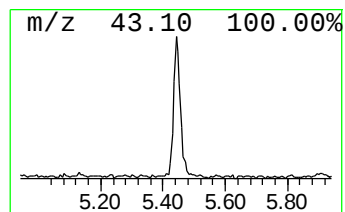
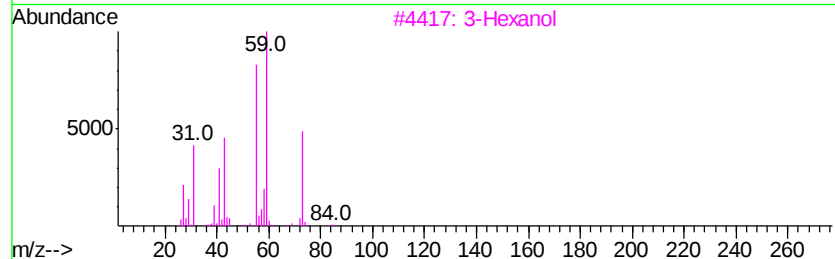
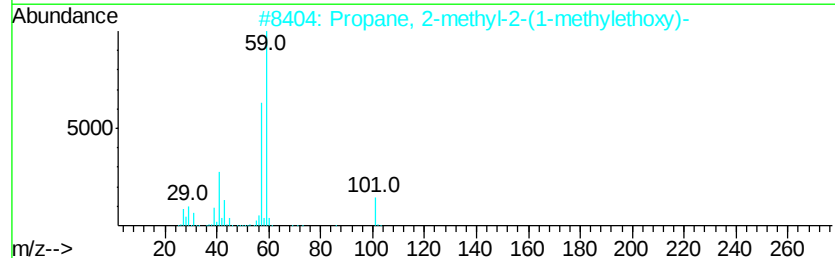
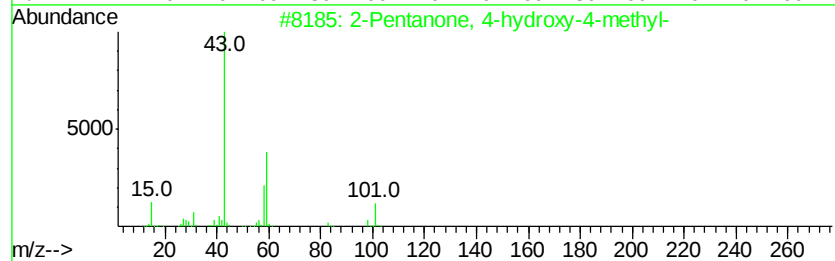
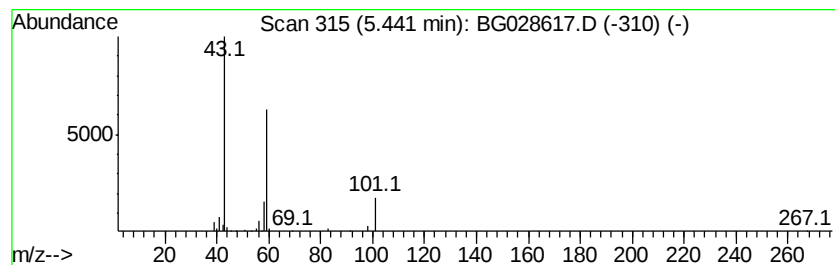
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TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.44	8.29 ng	79632	1,4-Dichlorobenzene-d4	8.33

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2		000123-42-2	53
2	Propane, 2-methyl-2-(1-methyleth...	116	C7H16O		017348-59-3	33
3	3-Hexanol	102	C6H14O		000623-37-0	33
4	3-Octanol	130	C8H18O		000589-98-0	33
5	2-Hexanol, 2-methyl-	116	C7H16O		000625-23-0	28



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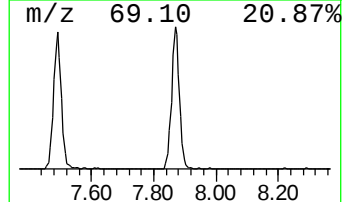
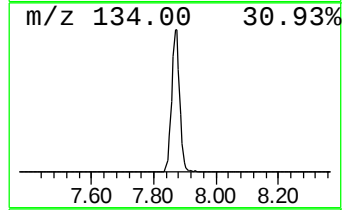
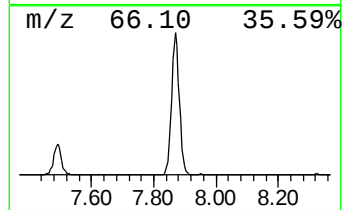
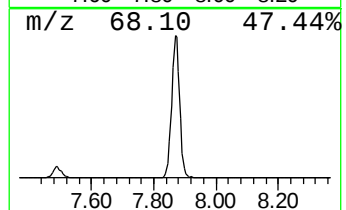
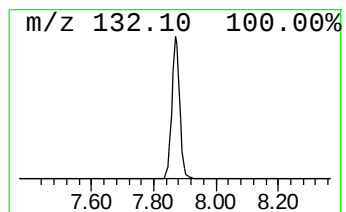
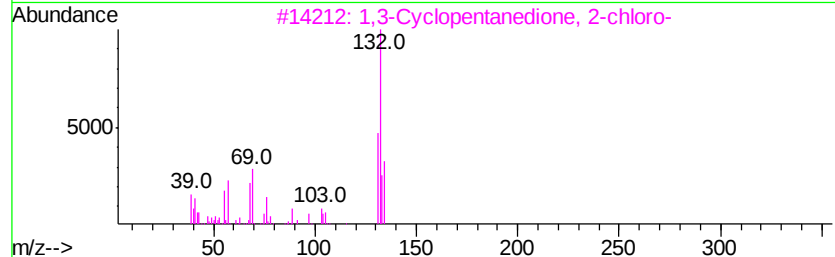
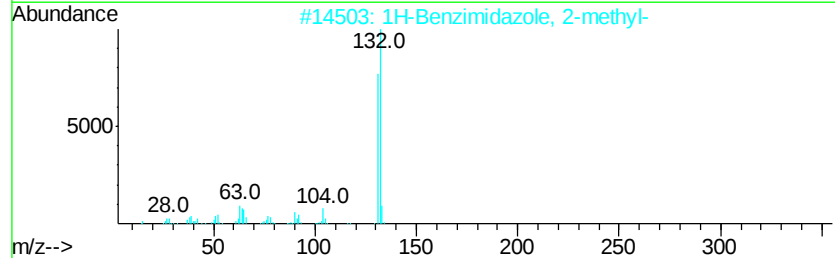
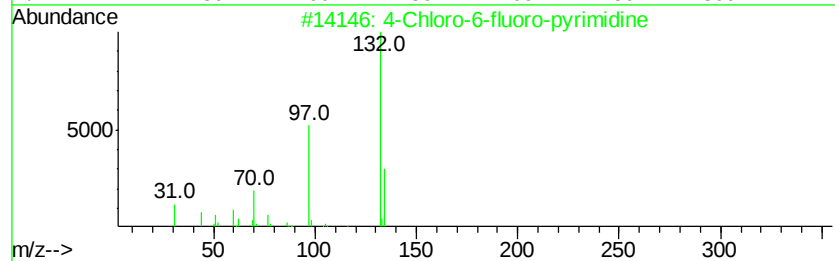
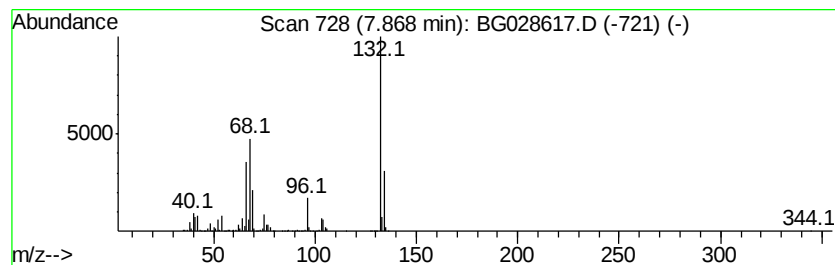
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Peak Number 2 unknown7.87 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.87	110.70 ng	1063690	1,4-Dichlorobenzene-d4	8.33

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	4-Chloro-6-fluoro-pyrimidine	132	C4H2ClFN2	051422-01-6	27
2		1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	25
3		1,3-Cyclopentanedione, 2-chloro-	132	C5H5ClO2	014203-19-1	25
4		(5-Methyl-2-pyridyl)acetonitrile	132	C8H8N2	1000241-93-9	10
5		1-Methylpyrrolo[1,2-a]pyrazine	132	C8H8N2	064608-59-9	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.44	8.3	ng	79632	1	8.33	192174	20.0
unknown7.87	7.87	110.7	ng	1063690	1	8.33	192174	20.0