

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG071718\
 Data File : BG035717.D
 Acq On : 18 Jul 2018 14:16
 Operator : JU/SJ
 Sample : PB111100TB
 Misc :
 ALS Vial : 36 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB111100TB

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-BG071218.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	5.146	310	316	324	rBV	111653	173298	3.57%	0.872%
2	5.622	388	397	412	rBV	879973	1457871	30.01%	7.335%
3	7.208	658	667	678	rBV	901698	1549883	31.91%	7.798%
4	7.573	719	729	742	rBV	883903	1511633	31.12%	7.605%
5	8.037	801	808	816	rBV2	153359	268687	5.53%	1.352%
6	8.360	856	863	872	rBV	662332	1149539	23.66%	5.783%
7	9.200	998	1006	1021	rBV	509746	940398	19.36%	4.731%
8	10.845	1278	1286	1293	rBV	172510	330483	6.80%	1.663%
9	13.282	1694	1701	1713	rBV	1446490	2306581	47.48%	11.605%
10	14.663	1929	1936	1948	rBV2	312128	511458	10.53%	2.573%
11	16.149	2182	2189	2208	rBV	1261836	1990293	40.97%	10.013%
12	17.406	2397	2403	2411	rBV2	507062	793064	16.33%	3.990%
13	20.015	2840	2847	2859	rBV	3698653	4857779	100.00%	24.440%
14	21.671	3122	3129	3140	rBV	607009	1057080	21.76%	5.318%
15	24.879	3664	3675	3688	rBV2	336556	978387	20.14%	4.922%

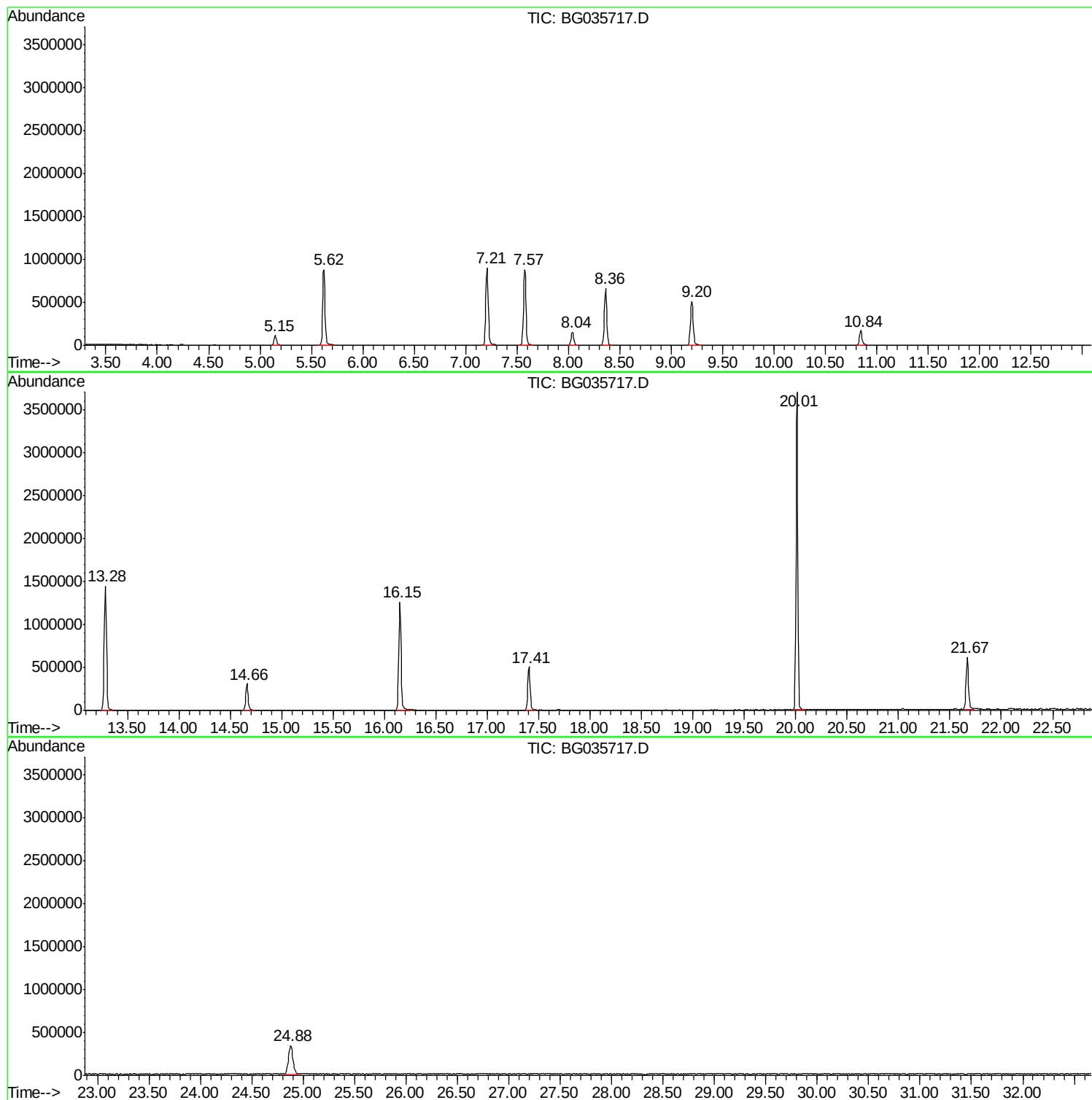
Sum of corrected areas: 19876434

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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P



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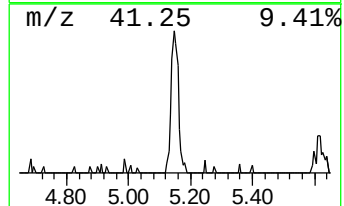
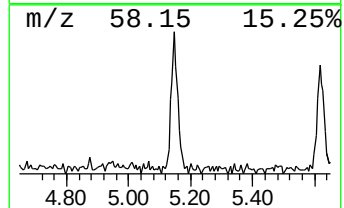
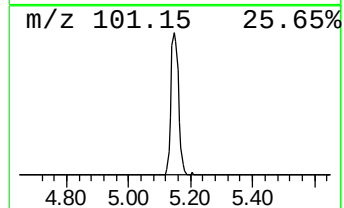
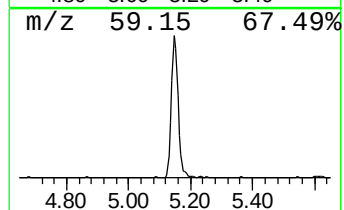
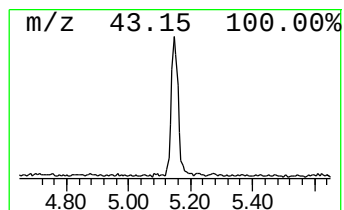
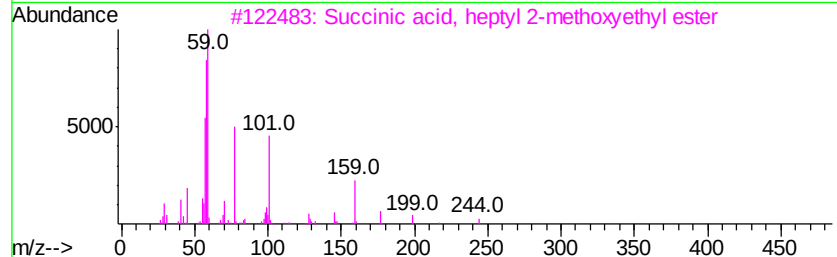
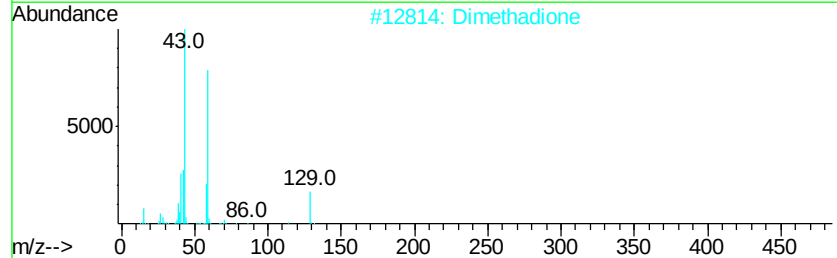
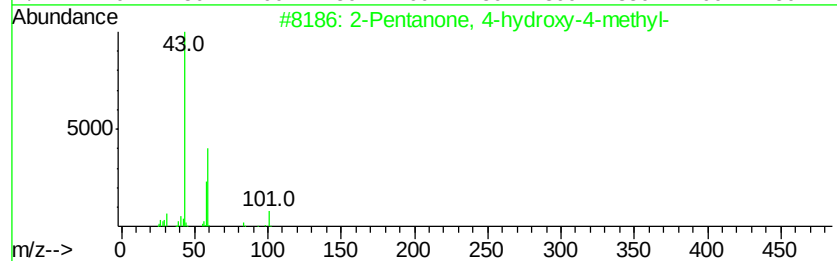
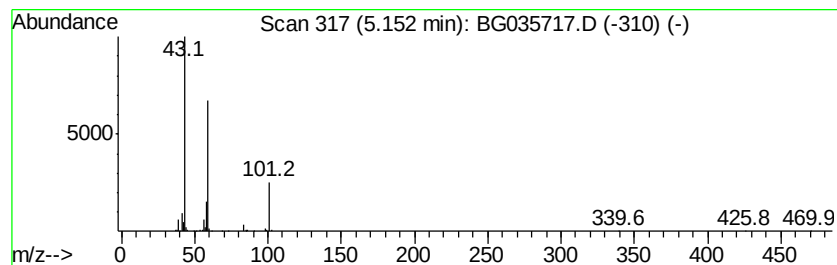
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.15	12.90 ng	173298	1,4-Dichlorobenzene-d4	8.04

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56
2		Dimethadione	129	C5H7NO3	000695-53-4	33
3		Succinic acid, heptyl 2-methoxyve...	274	C14H26O5	1000325-80-7	33
4		5-Hexen-2-one	98	C6H10O	000109-49-9	10
5		Butane	58	C4H10	000106-97-8	9



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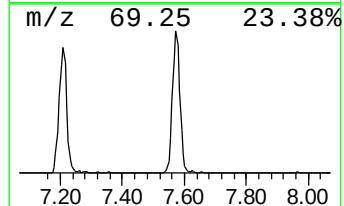
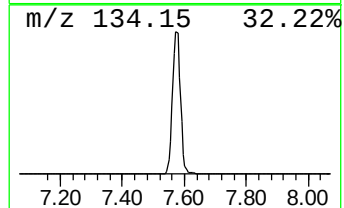
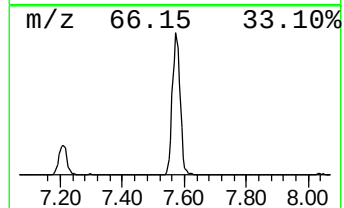
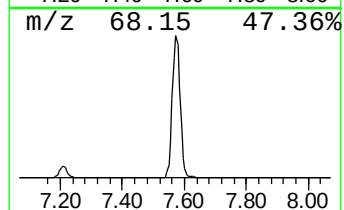
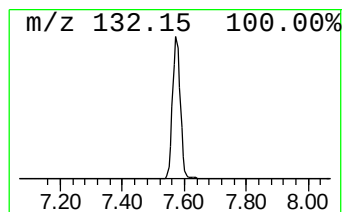
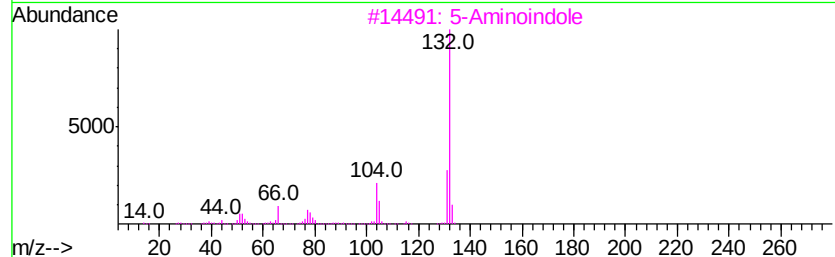
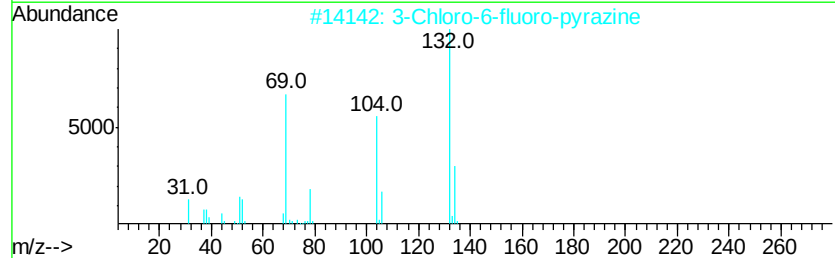
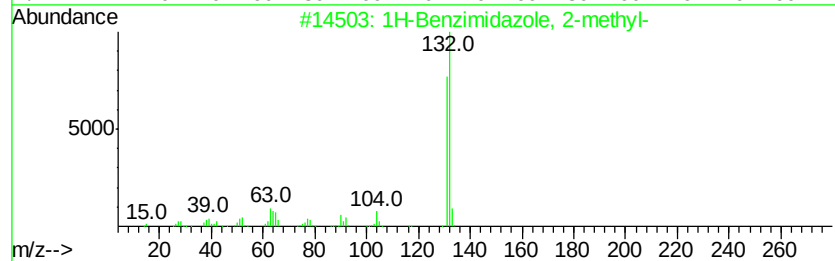
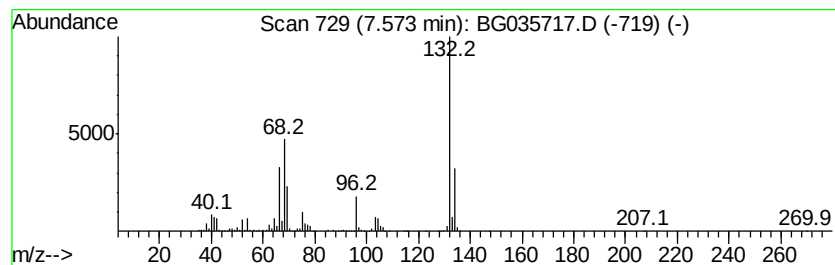
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Peak Number 2 unknown7.57 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.57	112.52 ng	1511630	1,4-Dichlorobenzene-d4	8.04

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	22
2		3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	16
3		5-Aminoindole	132	C8H8N2	005192-03-0	12
4		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
5		Cyclopropanamine, 1-phenyl-	133	C9H11N	1000351-26-2	10



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TIC Top Hit name	RT	EstConc	Units	Response	---Internal Standard---			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.15	12.9	ng	173298	1	8.04	268687	20.0
unknown7.57	7.57	112.5	ng	1511630	1	8.04	268687	20.0