

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG061818\
 Data File : BG035217.D
 Acq On : 18 Jun 2018 13:58
 Operator : SJ/JU
 Sample : J3504-06
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SB-51COMP

Integration Parameters: rteint.p
 Integrator: RTE
 Smoothing : ON
 Sampling : 1
 Start Thrs: 0.2
 Stop Thrs : 0

Filtering: 5
 Min Area: 3 % of largest Peak
 Max Peaks: 100
 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-BG060718.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	5.176	315	321	327	rBV	114820	169110	3.42%	0.783%
2	5.635	392	399	408	rBV	968458	1570438	31.74%	7.268%
3	7.221	660	669	677	rBV	854521	1480317	29.92%	6.851%
4	7.597	724	733	744	rVB	1032469	1769866	35.77%	8.191%
5	8.061	806	812	821	rVB	215472	354296	7.16%	1.640%
6	8.384	859	867	874	rBV	824177	1413531	28.57%	6.542%
7	9.224	1002	1010	1020	rVB	660453	1182886	23.90%	5.474%
8	10.863	1282	1289	1296	rBV	231883	430617	8.70%	1.993%
9	12.519	1565	1571	1577	rVB	32328	52252	1.06%	0.242%
10	13.307	1698	1705	1715	rBV	1677253	2581096	52.16%	11.945%
11	14.681	1933	1939	1947	rBV2	368425	586582	11.85%	2.715%
12	16.167	2184	2192	2211	rBV	1333921	2081928	42.07%	9.635%
13	17.424	2399	2406	2412	rBV	506119	758892	15.34%	3.512%
14	20.033	2844	2850	2856	rBV	3791988	4948371	100.00%	22.900%
15	21.689	3125	3132	3140	rBV	722030	1167438	23.59%	5.403%
16	24.920	3671	3682	3694	rBV	370068	1060706	21.44%	4.909%

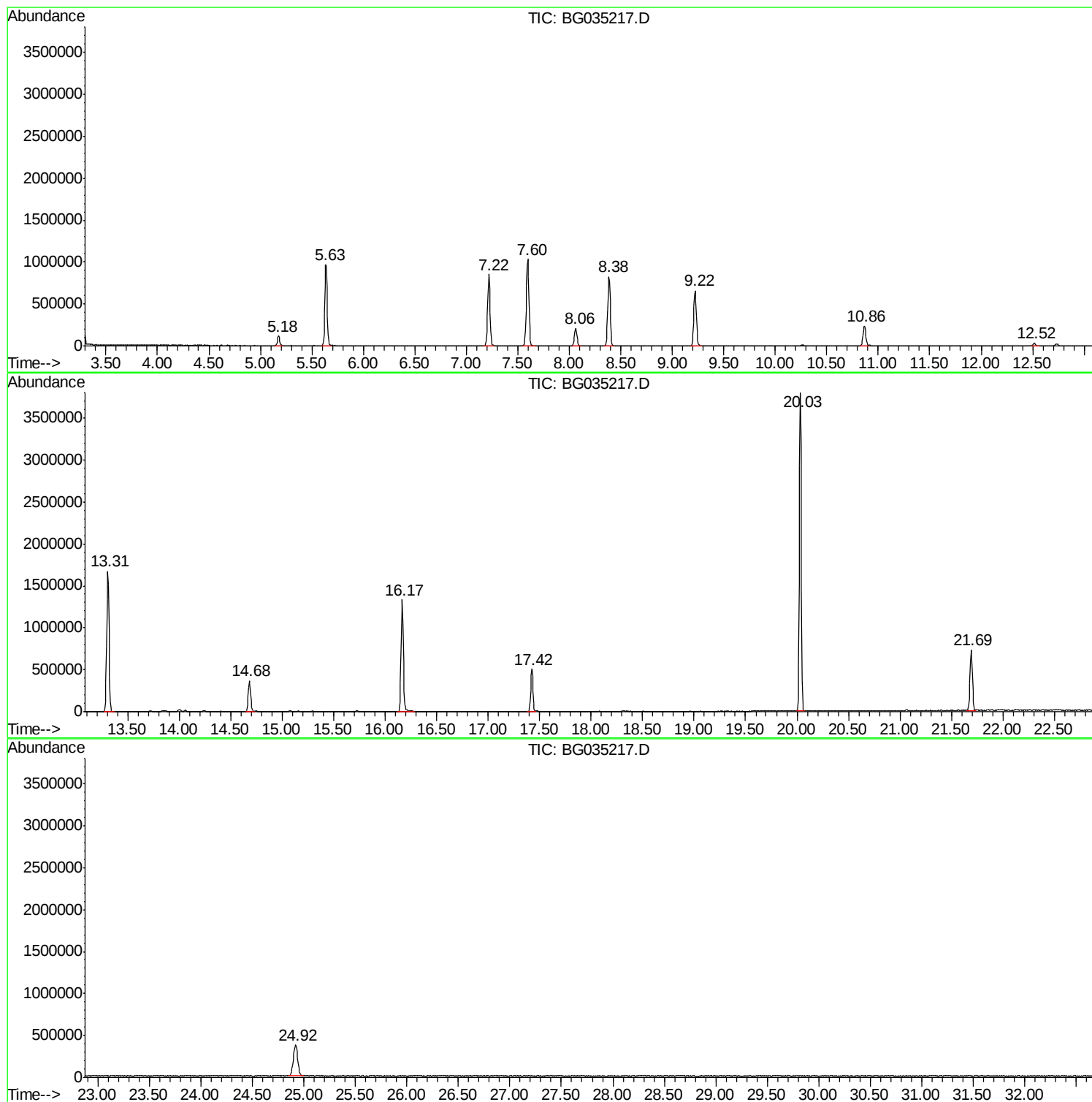
Sum of corrected areas: 21608326

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Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG060718.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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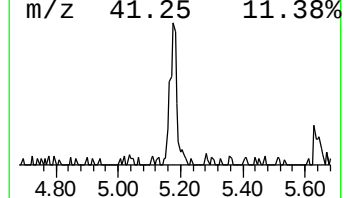
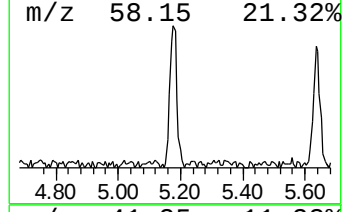
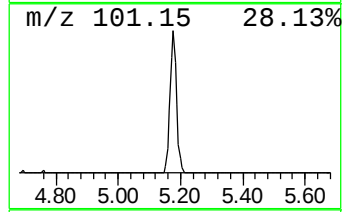
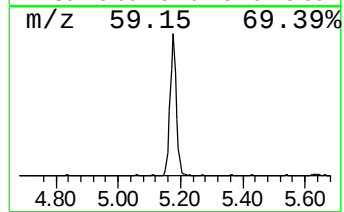
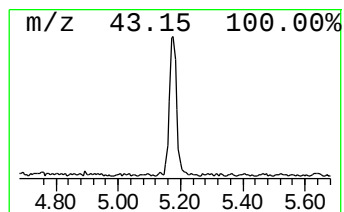
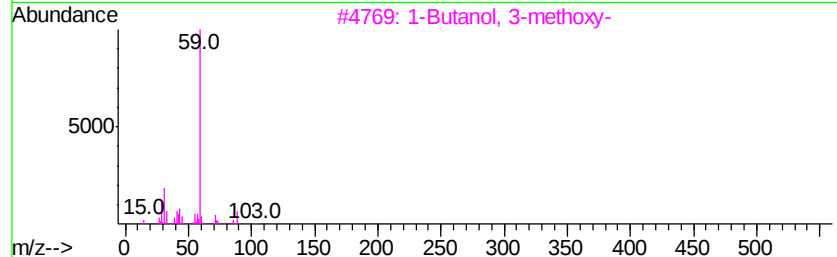
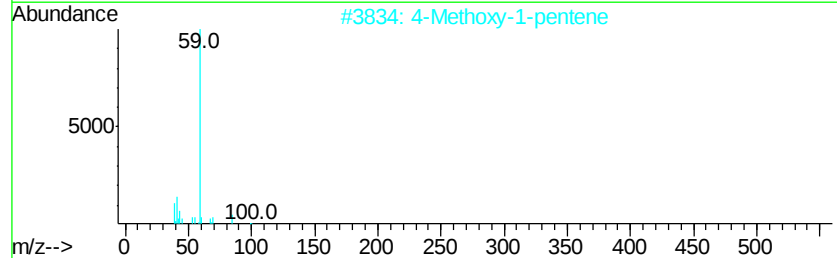
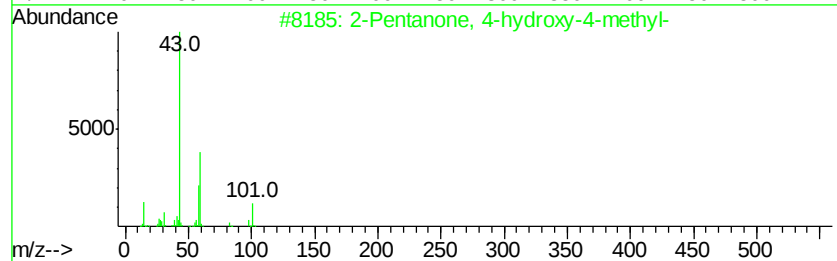
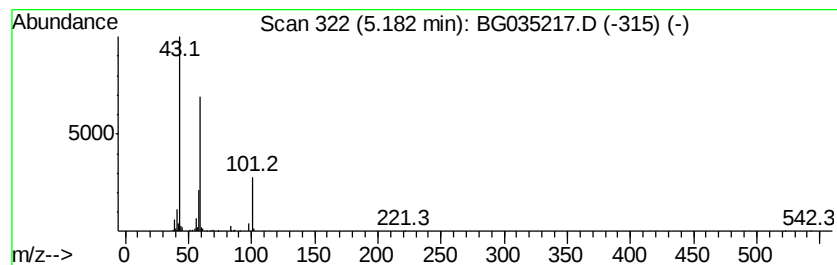
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.18	9.55 ng	169110	1,4-Dichlorobenzene-d4	8.07

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	72
2			4-Methoxy-1-pentene	100	C6H12O	098386-09-5	43
3			1-Butanol, 3-methoxy-	104	C5H12O2	002517-43-3	37
4			Hydrazine, 1,1-bis(1-methylethyl)-	116	C6H16N2	000921-14-2	36
5			2,5,8,11-Tetraoxadodecane	178	C8H18O4	000112-49-2	33



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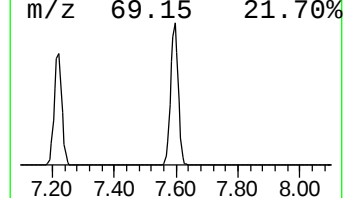
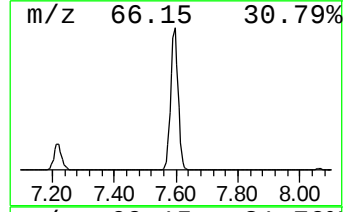
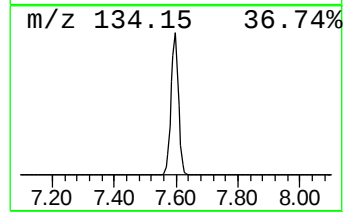
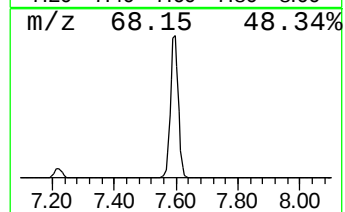
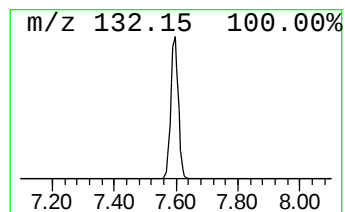
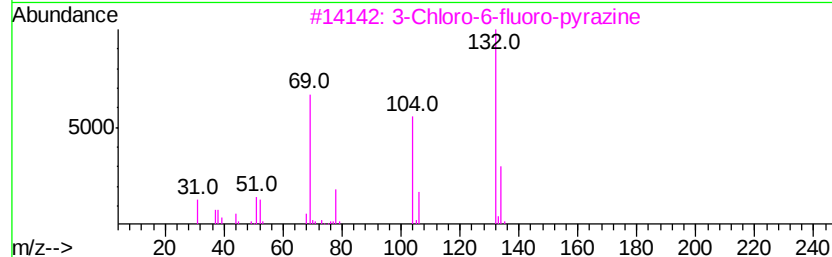
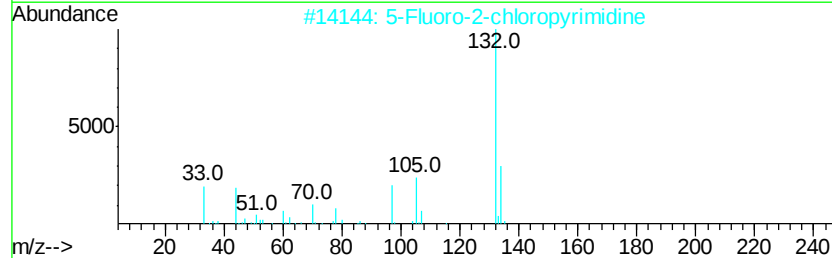
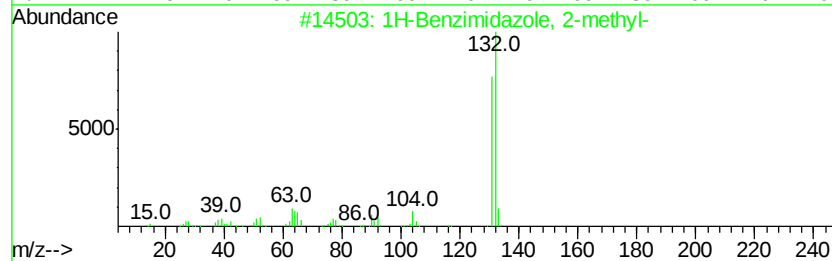
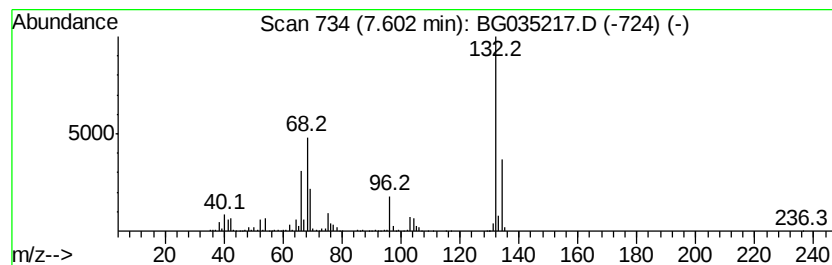
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Peak Number 2 unknown7.60 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.60	99.91 ng	1769870	1,4-Dichlorobenzene-d4	8.07

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	22
2		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	16
3		3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	16
4		5-Aminoindole	132	C8H8N2	005192-03-0	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...	5.18	9.6	ng	169110	1	8.07	354296	20.0
unknown7.60	7.60	99.9	ng	1769870	1	8.07	354296	20.0