

Data Path : Z:\HPCHEM1\BNA G\DATA\BG042018\
 Data File : BG033887.D
 Acq On : 20 Apr 2018 15:40
 Operator : SJ/JU
 Sample : J2460-04
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 MDP-B-2-18-20

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-BG041918.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	5.000	285	291	296	rBV	46491	72763	2.65%	0.389%
2	5.446	358	367	375	rBV	924115	1428979	52.05%	7.647%
3	7.003	624	632	644	rBV	886638	1532597	55.82%	8.201%
4	7.368	686	694	702	rBV	888035	1486007	54.12%	7.952%
5	7.838	765	774	781	rBV	256975	445508	16.23%	2.384%
6	8.155	818	828	836	rBV	577032	1002212	36.50%	5.363%
7	8.983	961	969	978	rBV	495531	868132	31.62%	4.646%
8	10.617	1239	1247	1255	rVB	360274	632919	23.05%	3.387%
9	13.090	1661	1668	1676	rVB	1121869	1702817	62.02%	9.112%
10	13.925	1805	1810	1816	rBV	79738	123469	4.50%	0.661%
11	14.389	1884	1889	1894	rBV2	21288	30119	1.10%	0.161%
12	14.465	1896	1902	1911	rVB2	555382	880003	32.05%	4.709%
13	15.346	2048	2052	2056	rBV2	22901	30244	1.10%	0.162%
14	15.957	2149	2156	2165	rBV	1091216	1652812	60.20%	8.844%
15	17.215	2363	2370	2377	rBV2	747415	1139929	41.52%	6.100%
16	18.137	2522	2527	2534	rBV3	45460	79711	2.90%	0.427%
17	19.565	2765	2770	2776	rVB	26077	35486	1.29%	0.190%
18	19.853	2813	2819	2826	rVB	2099301	2745553	100.00%	14.692%
19	21.474	3088	3095	3103	rVB2	907474	1393098	50.74%	7.455%
20	23.173	3379	3384	3391	rVB	22832	45489	1.66%	0.243%
21	24.542	3606	3617	3630	rVB2	499325	1359739	49.53%	7.276%

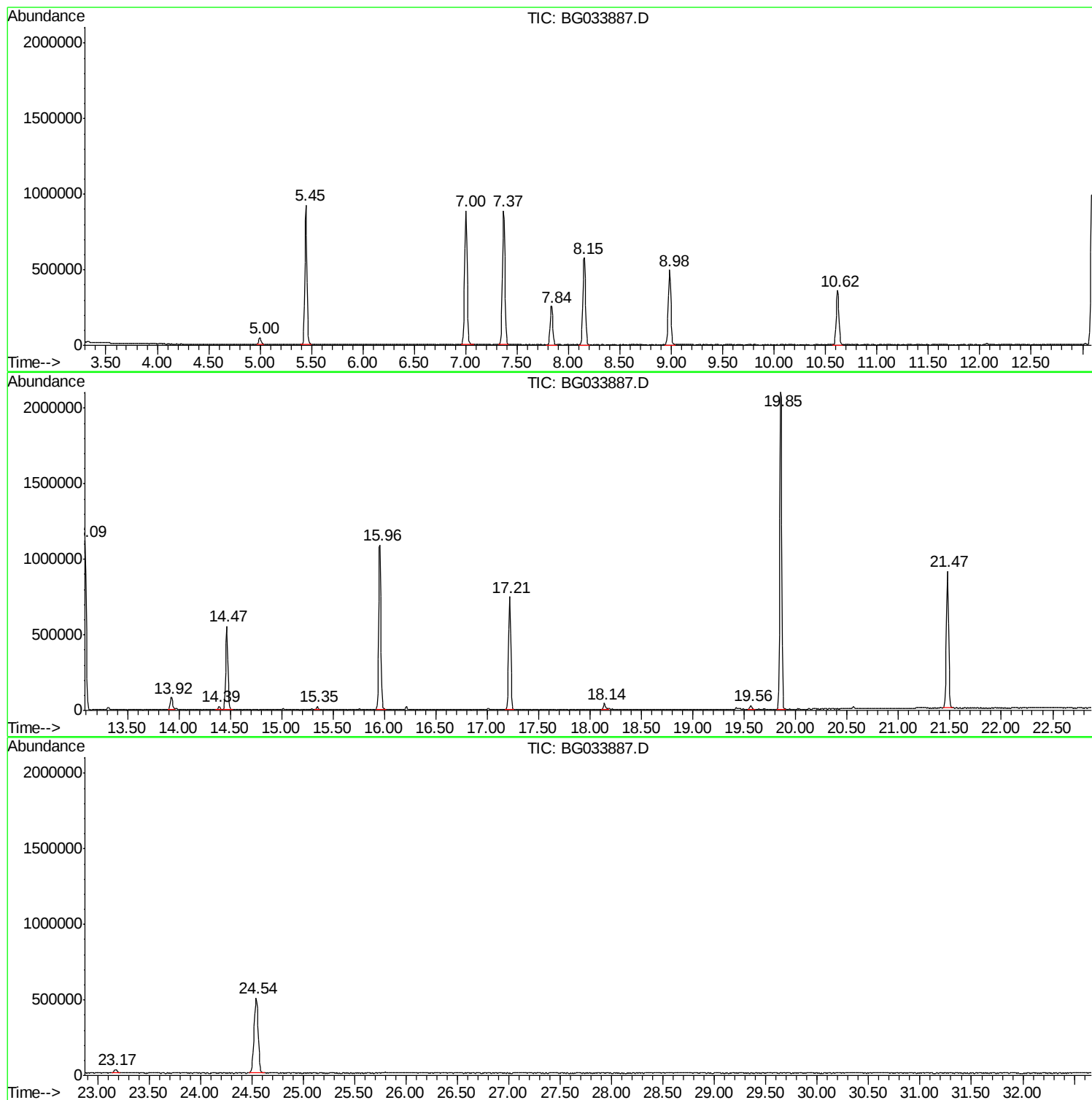
Sum of corrected areas: 18687586

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Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG041918.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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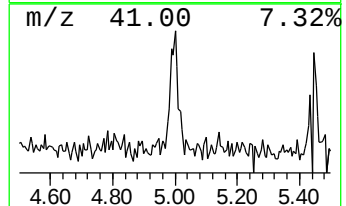
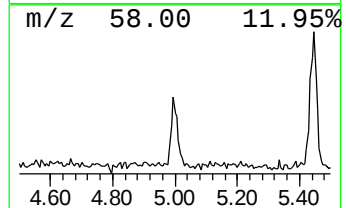
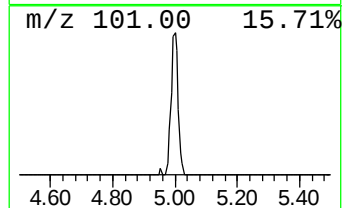
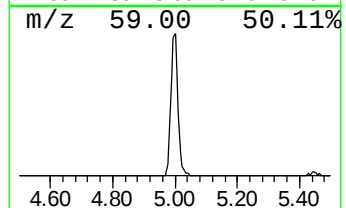
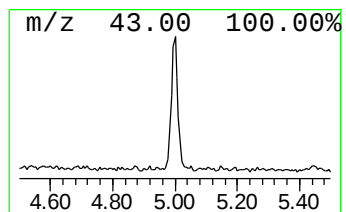
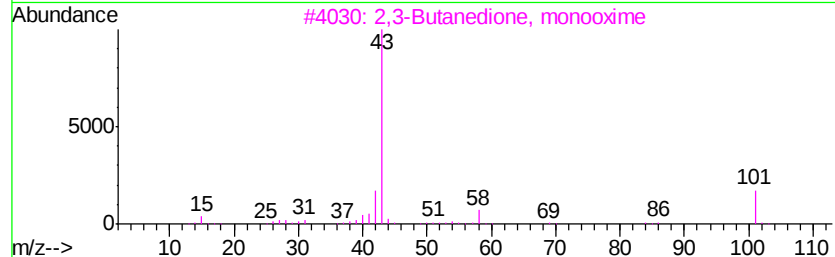
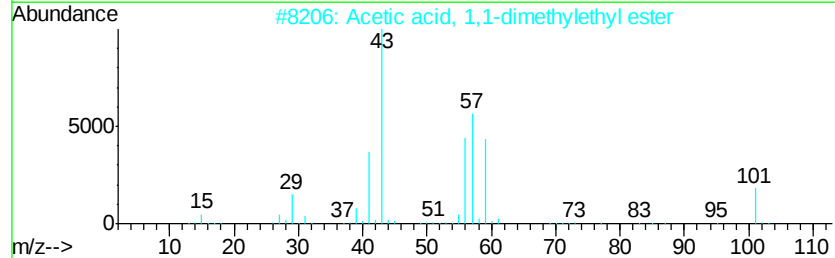
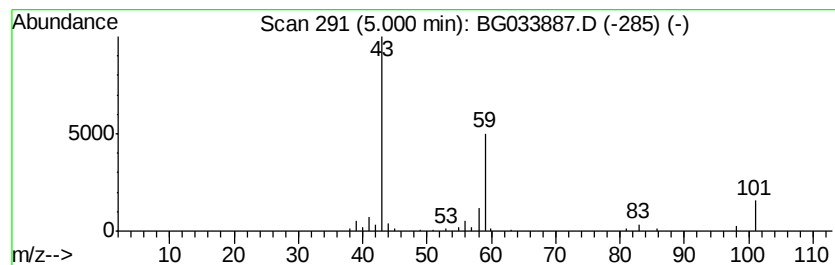
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.00	3.27 ng	72763	1,4-Dichlorobenzene-d4	7.83

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	64
2			Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	50
3			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	35
4			5-Hexen-2-one	98	C6H10O	000109-49-9	9
5			3-Pentanol	88	C5H12O	000584-02-1	9



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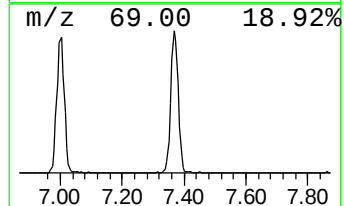
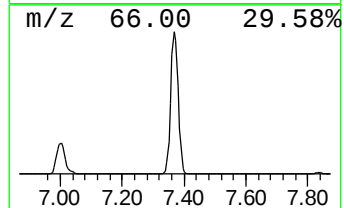
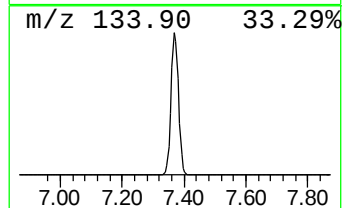
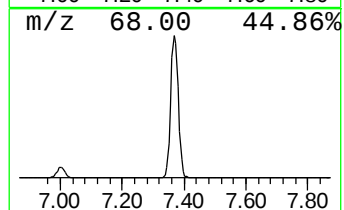
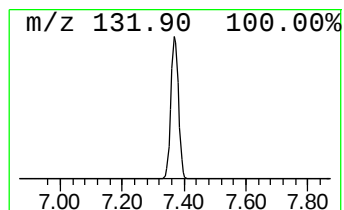
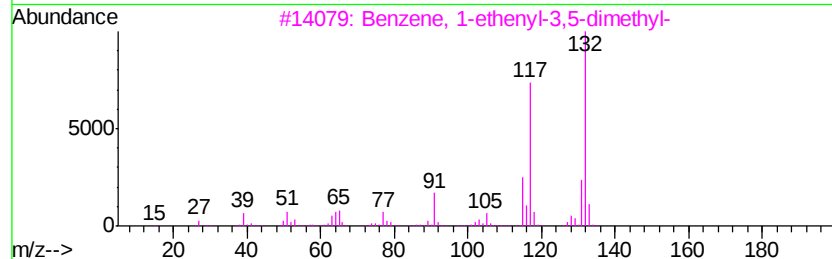
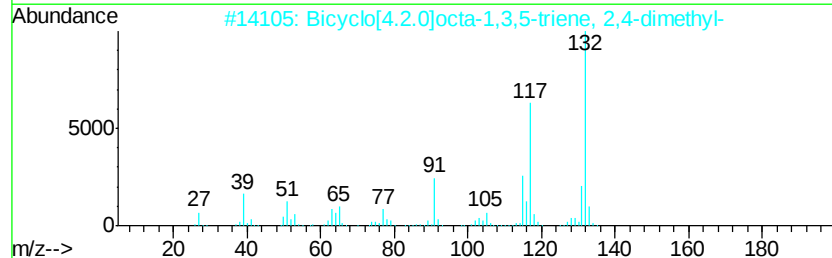
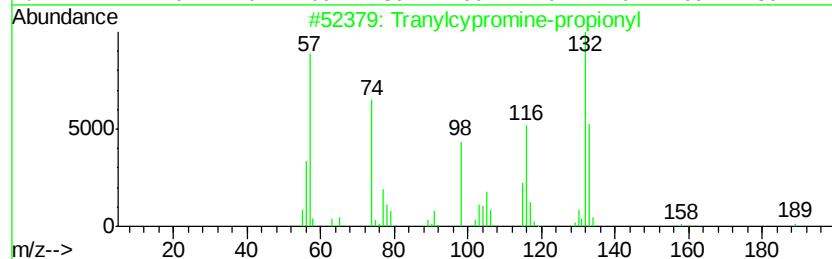
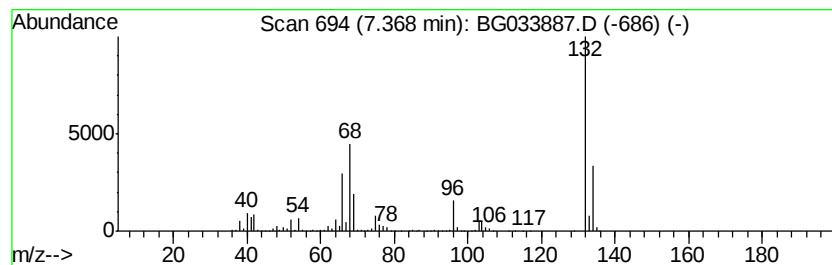
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Peak Number 2 unknown7.37 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.37	66.71 ng	1486010	1,4-Dichlorobenzene-d4	7.83

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tranvlcvpromine-propionyl	189	C12H15NO	1000123-86-3	10
2		Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9
3		Benzene, 1-ethenyl-3,5-dimethyl-	132	C10H12	005379-20-4	9
4		Xenon	132	Xe	007440-63-3	9
5		4-Chloro-2-fluoro-pyrimidine	132	C4H2ClFN2	051422-00-5	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.00	3.3	ng	72763	1	7.83	445508	20.0
unknown7.37	7.37	66.7	ng	1486010	1	7.83	445508	20.0