

Data Path : Z:\HPCHEM1\BNA G\DATA\BG042418\
 Data File : BG033962.D
 Acq On : 24 Apr 2018 20:04
 Operator : SJ/JU
 Sample : J2475-03
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 MW-02

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	4.991	285	290	299	rBV	69502	105094	1.95%	0.392%
2	5.437	360	366	379	rBV	972038	1532487	28.42%	5.713%
3	6.994	623	631	642	rBV	691905	1145238	21.24%	4.270%
4	7.364	685	694	704	rBV	1354354	2299028	42.63%	8.571%
5	7.829	765	773	780	rBV	257256	437195	8.11%	1.630%
6	8.146	819	827	836	rBV	1151732	1946024	36.08%	7.255%
7	8.974	958	968	977	rBV	874830	1545916	28.67%	5.763%
8	10.608	1237	1246	1255	rVB	358471	618951	11.48%	2.307%
9	13.081	1659	1667	1675	rBV	2547358	3782580	70.14%	14.102%
10	14.456	1895	1901	1911	rBV2	550642	865793	16.05%	3.228%
11	15.949	2148	2155	2177	rBV	1945910	2948285	54.67%	10.991%
12	17.212	2363	2370	2379	rBV	719383	1130440	20.96%	4.214%
13	18.122	2516	2525	2534	rBV6	44150	88556	1.64%	0.330%
14	19.844	2812	2818	2825	rBV	4051583	5392965	100.00%	20.105%
15	21.466	3087	3094	3104	rVB	943000	1534291	28.45%	5.720%
16	24.521	3603	3614	3628	rVB2	536557	1450719	26.90%	5.408%

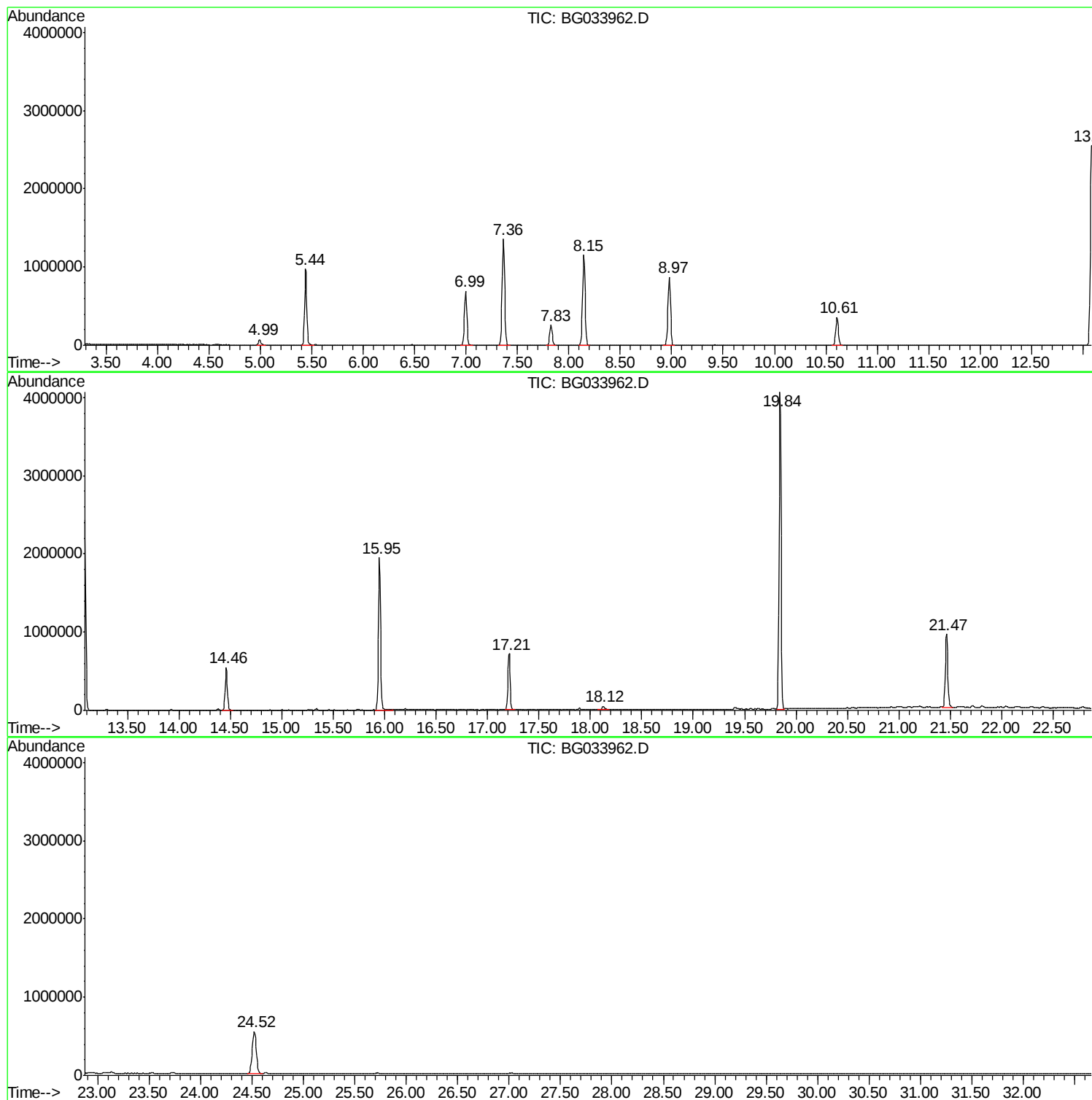
Sum of corrected areas: 26823562

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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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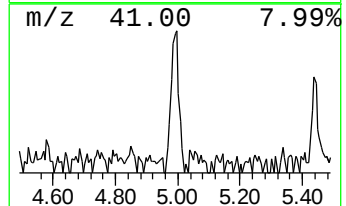
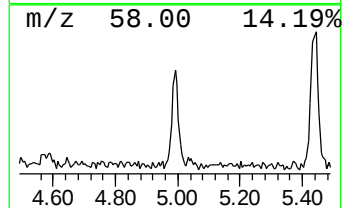
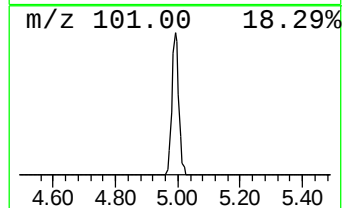
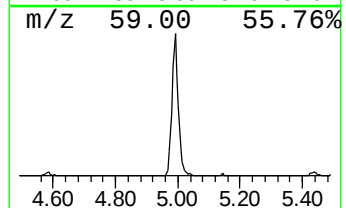
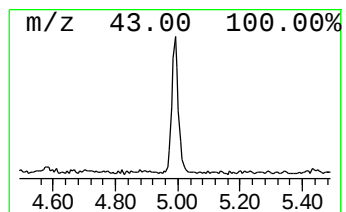
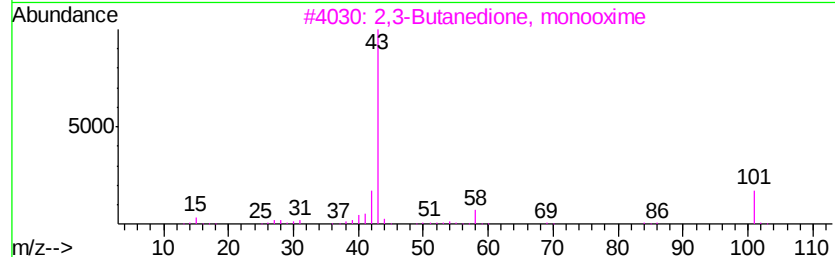
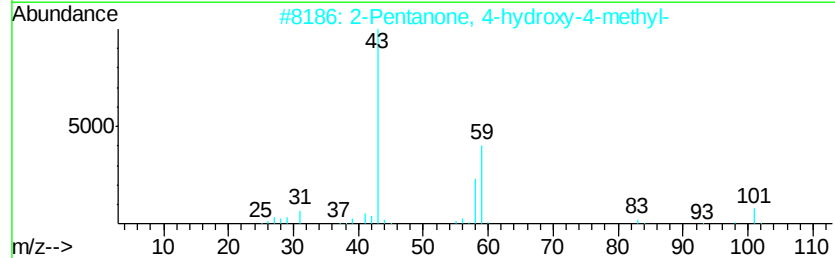
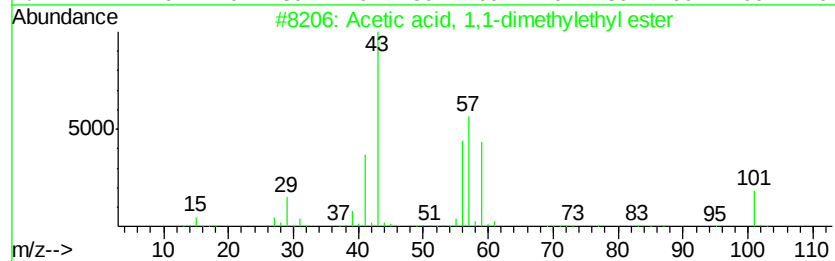
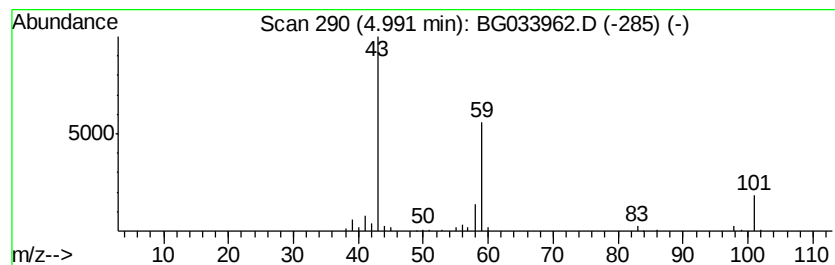
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Peak Number 1 Acetic acid, 1,1-dimethylet... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.99	4.81 ng	105094	1,4-Dichlorobenzene-d4	7.83

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	50
2		2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	45
3		2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	27
4		3-Heptanol	116	C7H16O	000589-82-2	25
5		Propane, 2-ethoxy-2-methyl-	102	C6H14O	000637-92-3	23



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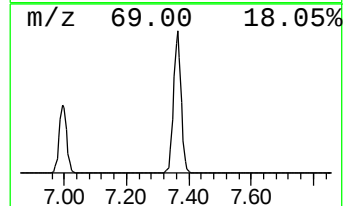
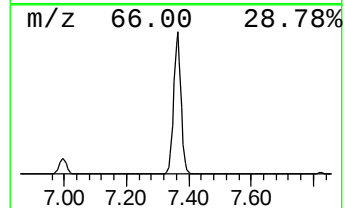
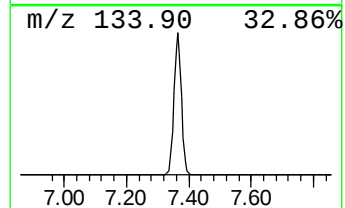
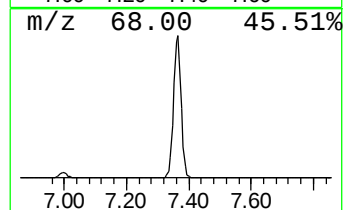
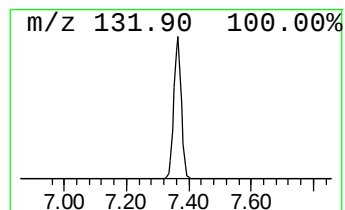
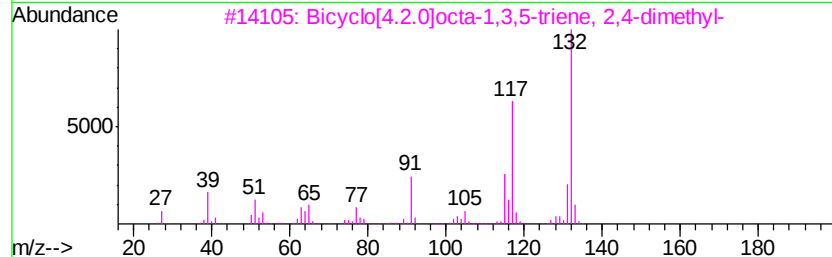
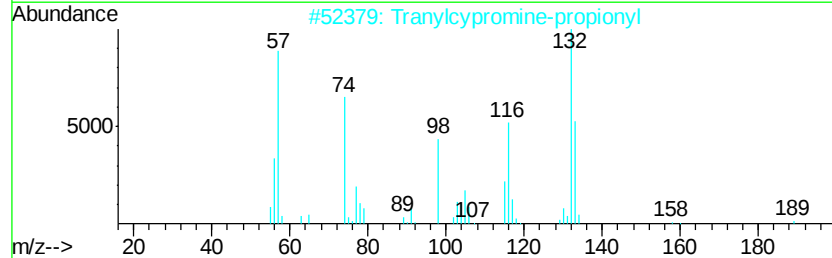
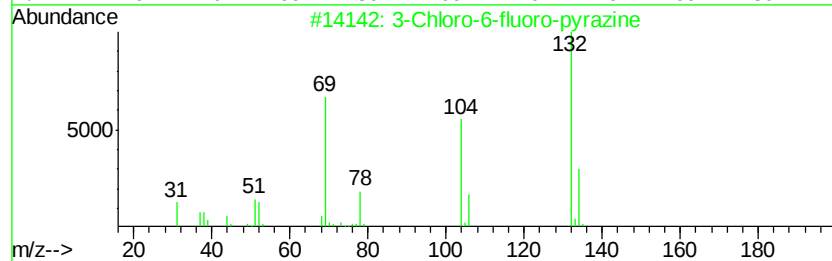
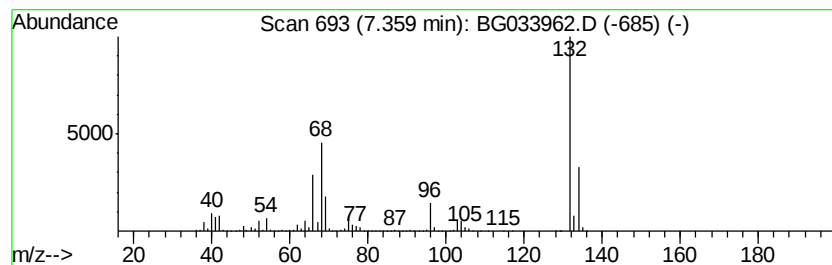
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Peak Number 2 unknown7.36 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.36	105.17 ng	2299030	1,4-Dichlorobenzene-d4	7.83

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
3		Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9
4		Benzene, 1-ethenyl-3,5-dimethyl-	132	C10H12	005379-20-4	9
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
Acetic acid, 1,1-...	4.99	4.8	ng	105094	1	7.83	437195	20.0
unknown7.36	7.36	105.2	ng	2299030	1	7.83	437195	20.0