

Data Path : Z:\HPCHEM1\BNA G\DATA\BG042018\
 Data File : BG033908.D
 Acq On : 21 Apr 2018 6:20
 Operator : SJ/JU
 Sample : J2473-11
 Misc :
 ALS Vial : 27 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 041818-SIDW-F-D-B

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	4.992	283	290	297	rBV	60360	90336	1.43%	0.411%
2	5.439	359	366	377	rBV	422695	663042	10.48%	3.017%
3	6.996	623	631	640	rBV	265486	449610	7.10%	2.046%
4	7.366	686	694	703	rBV	689903	1184254	18.71%	5.389%
5	7.830	766	773	780	rVB	216618	359554	5.68%	1.636%
6	8.148	818	827	837	rVB	840546	1414864	22.36%	6.439%
7	8.982	958	969	978	rBV	711069	1271273	20.09%	5.785%
8	10.615	1239	1247	1256	rBV	295213	521272	8.24%	2.372%
9	13.089	1661	1668	1675	rBV	1924861	3006556	47.51%	13.682%
10	14.464	1895	1902	1913	rVB2	495264	788099	12.45%	3.587%
11	15.950	2148	2155	2166	rBV	1291631	1865939	29.49%	8.492%
12	17.213	2363	2370	2377	rBV2	712158	1101647	17.41%	5.013%
13	18.130	2520	2526	2536	rBV2	46280	80538	1.27%	0.367%
14	19.857	2813	2820	2825	rBV	4651838	6328149	100.00%	28.798%
15	21.473	3087	3095	3104	rBV	958094	1448825	22.89%	6.593%
16	24.534	3605	3616	3628	rBV2	523664	1400023	22.12%	6.371%

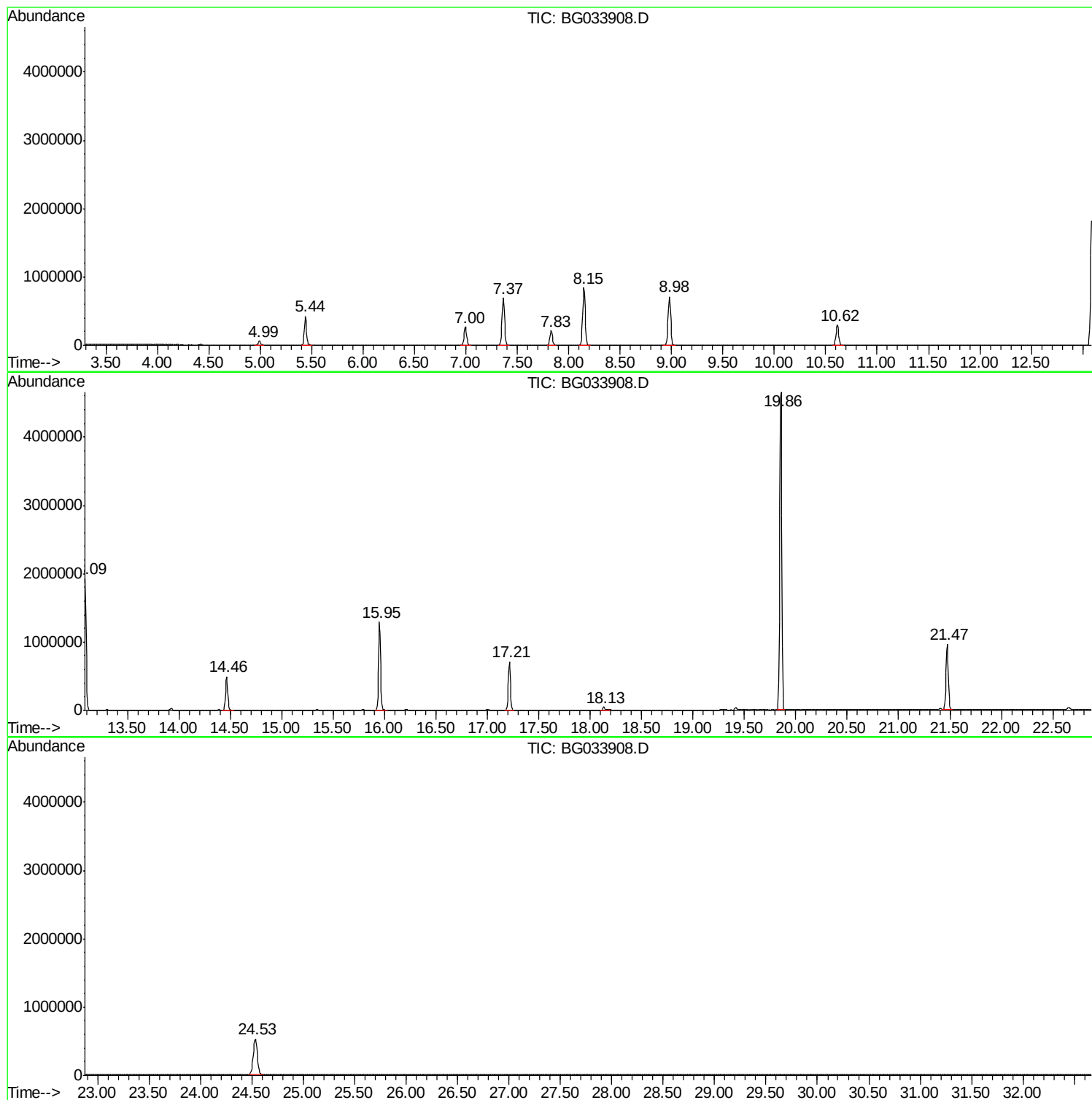
Sum of corrected areas: 21973981

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Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG041918.M
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TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P



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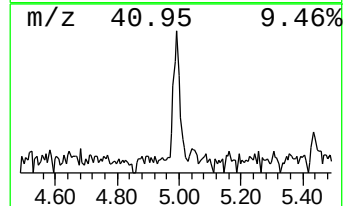
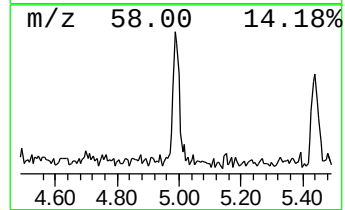
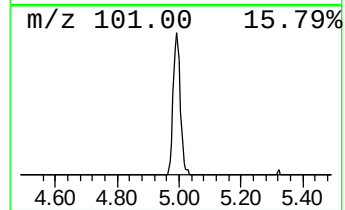
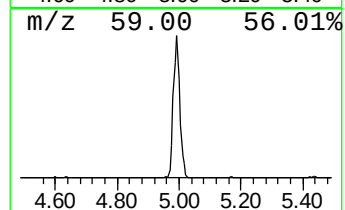
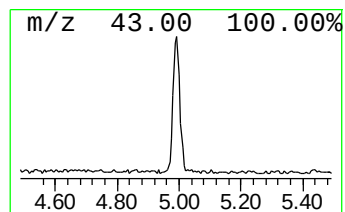
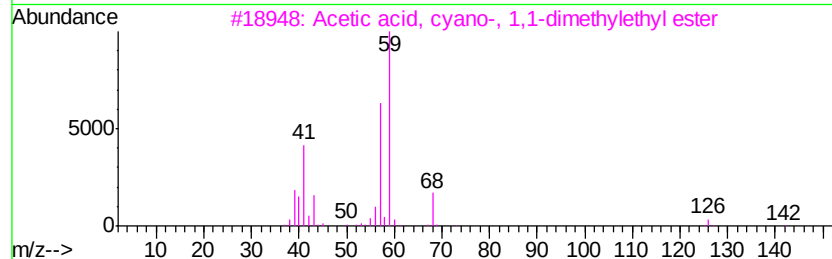
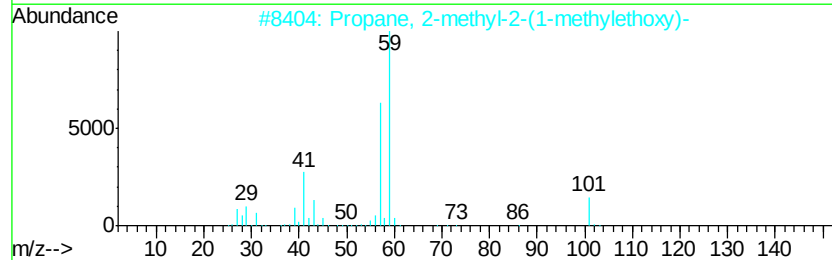
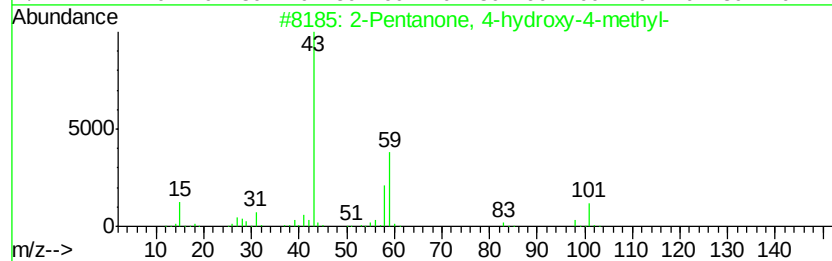
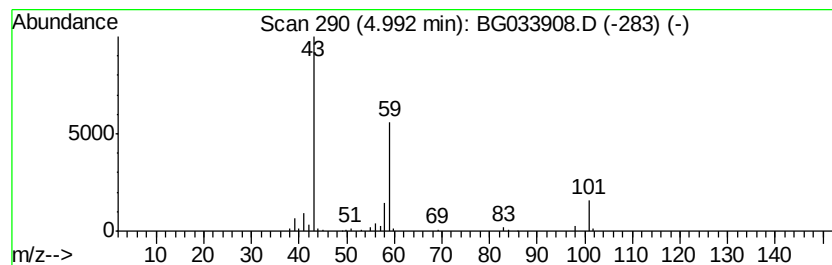
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.99	5.02 ng	90336	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	50	
2	Propane, 2-methyl-2-(1-methyleth...	116	C7H16O	017348-59-3	33	
3	Acetic acid, cyano-, 1,1-dimethv...	141	C7H11NO2	001116-98-9	23	
4	2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	16	
5	5-Hexen-2-one	98	C6H10O	000109-49-9	9	



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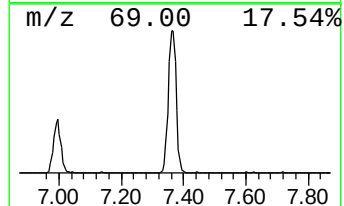
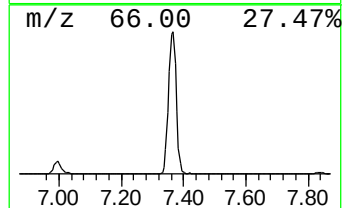
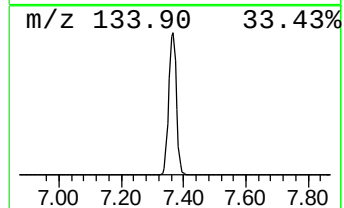
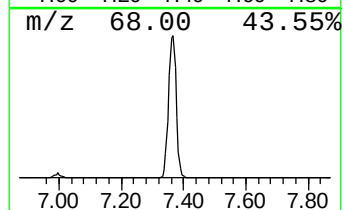
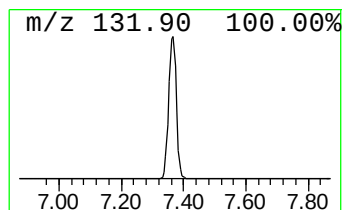
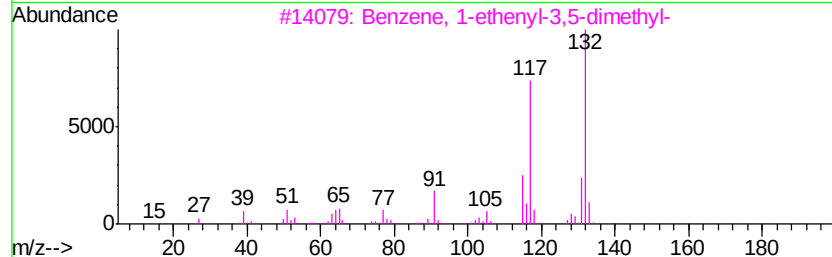
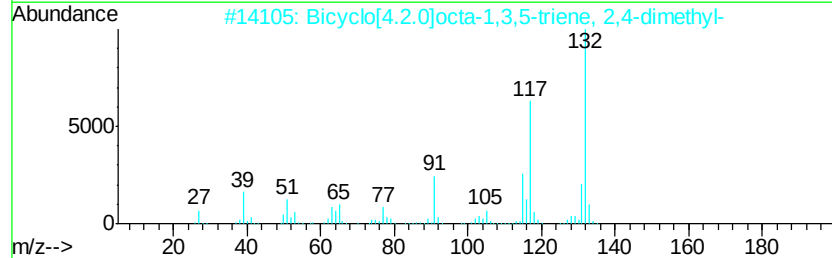
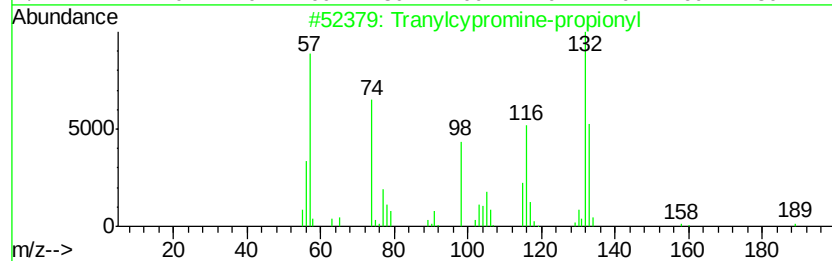
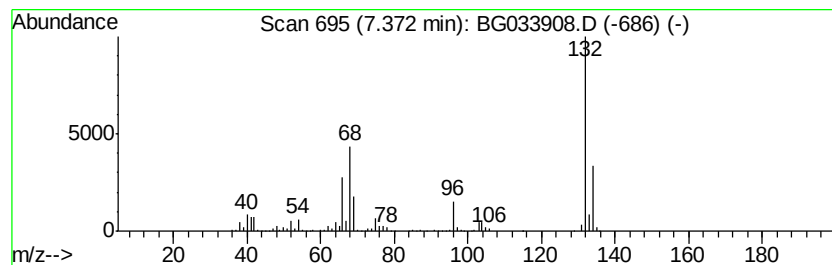
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Peak Number 2 unknown7.37 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.37	65.87 ng	1184250	1,4-Dichlorobenzene-d4	7.83

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tranlycypromine-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		2H-Cyclopenta[d]pyridazine, 2-me...	132	C8H8N2	022291-85-6	9
5		1,3,4-Thiadiazole-2(3H)-thione, ...	132	C3H4N2S2	029490-19-5	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	4.99	5.0	ng	90336	1	7.83	359554	20.0
unknown7.37	7.37	65.9	ng	1184250	1	7.83	359554	20.0