

Data Path : Z:\HPCHEM1\BNA G\DATA\BG042018\
 Data File : BG033906.D
 Acq On : 21 Apr 2018 5:05
 Operator : SJ/JU
 Sample : J2473-09
 Misc :
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 041818-SIDI-F-U-M

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	281	290	297	rVB	61620	96736	1.39%	0.407%
2	5.439	360	366	378	rVB	456505	698201	10.05%	2.938%
3	6.996	624	631	640	rVB	294617	482666	6.95%	2.031%
4	7.366	686	694	706	rVB	765164	1284746	18.49%	5.407%
5	7.830	765	773	781	rVB	218169	366814	5.28%	1.544%
6	8.147	819	827	837	rBV	898987	1550871	22.32%	6.527%
7	8.981	961	969	979	rBV	799865	1397107	20.11%	5.880%
8	10.615	1240	1247	1255	rBV	308580	535959	7.71%	2.256%
9	13.088	1660	1668	1678	rVB	2195278	3342172	48.10%	14.065%
10	14.463	1895	1902	1909	rBV2	517646	827062	11.90%	3.481%
11	15.950	2148	2155	2168	rBV	1374260	2064431	29.71%	8.688%
12	17.213	2364	2370	2378	rBV2	790480	1149951	16.55%	4.840%
13	18.130	2521	2526	2534	rBV3	46184	81384	1.17%	0.343%
14	19.857	2813	2820	2825	rBV	5259163	6948308	100.00%	29.242%
15	21.473	3088	3095	3102	rVB	957974	1480569	21.31%	6.231%
16	24.534	3605	3616	3629	rVB2	547443	1454627	20.93%	6.122%

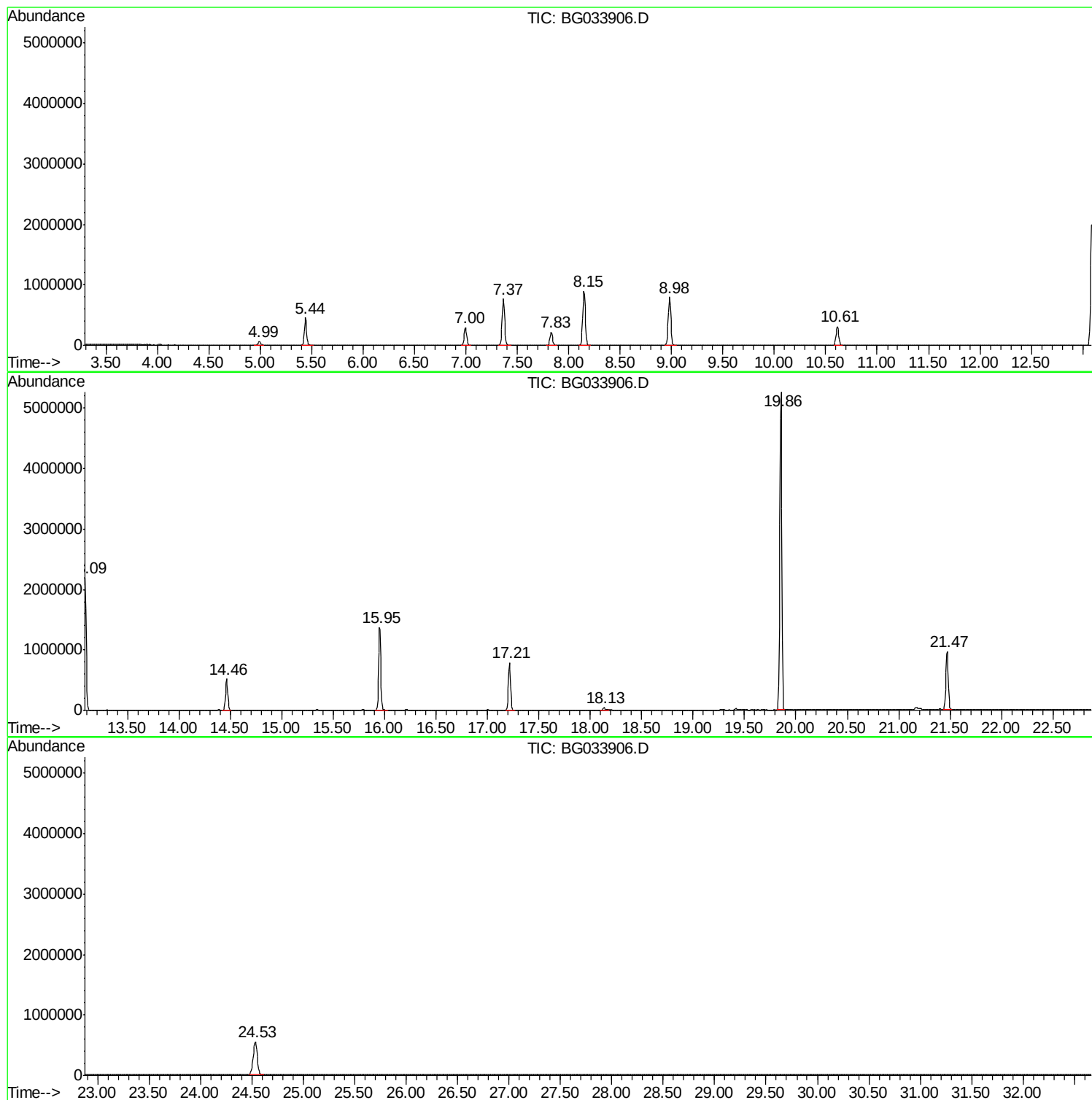
Sum of corrected areas: 23761604

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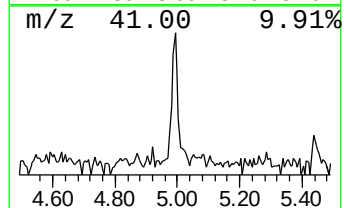
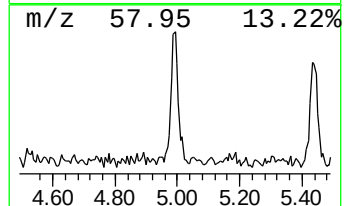
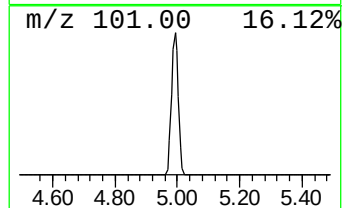
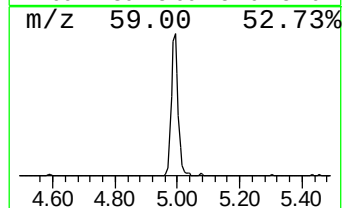
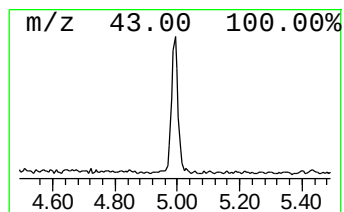
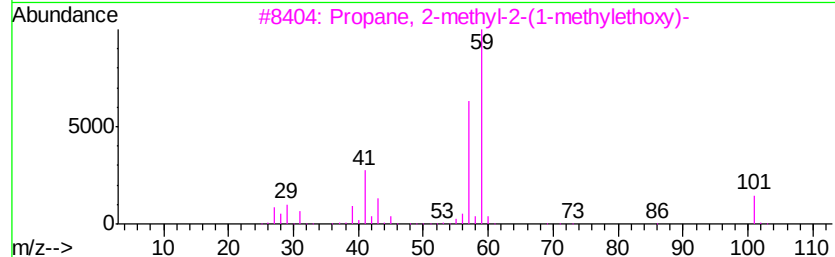
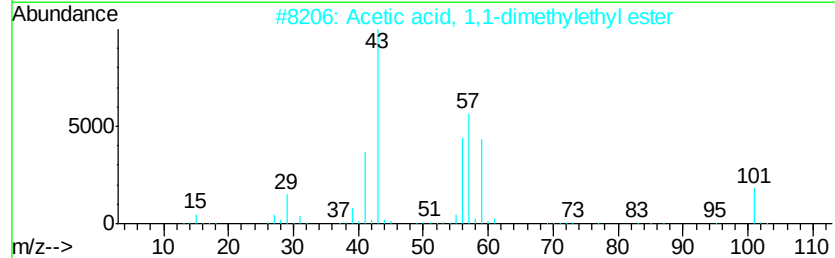
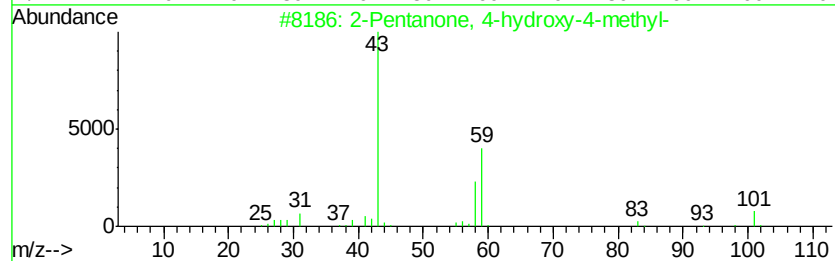
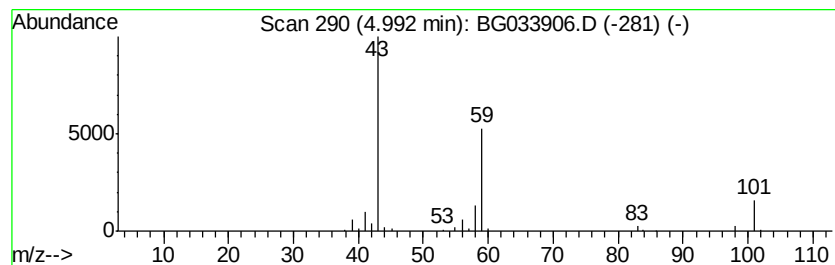
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.99	5.27 ng	96736	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			Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	38
3			Propane, 2-methyl-2-(1-methyleth...	116	C7H16O	017348-59-3	36
4			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9
5			5-Hexen-2-one	98	C6H10O	000109-49-9	9



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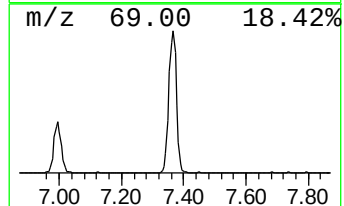
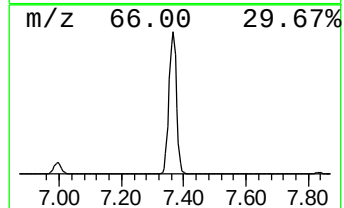
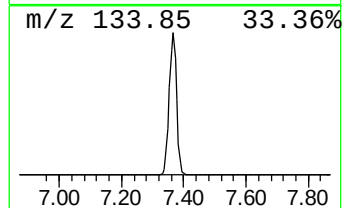
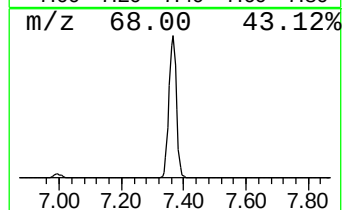
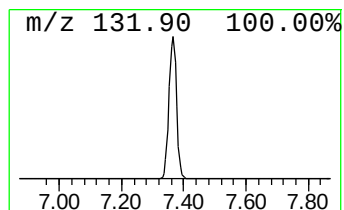
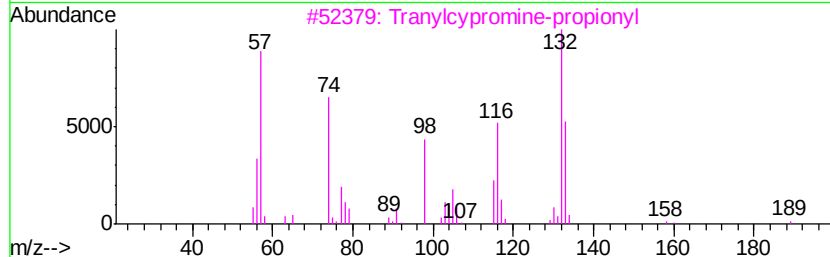
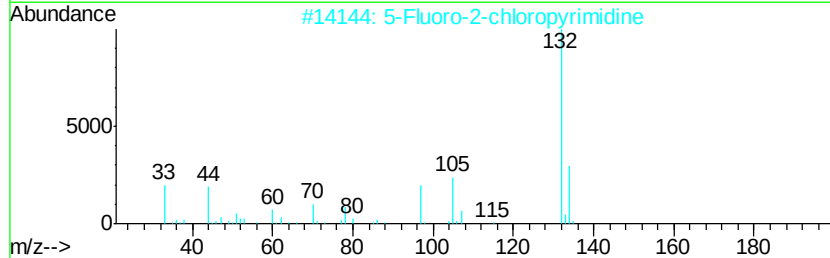
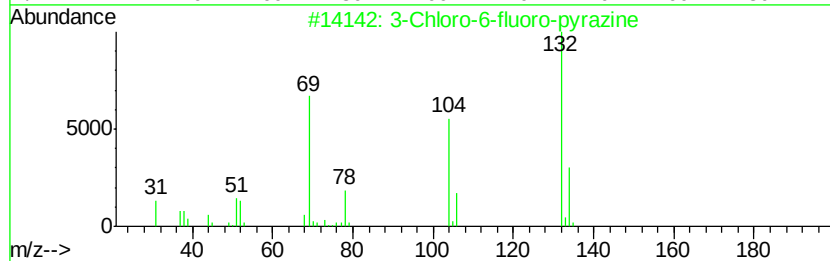
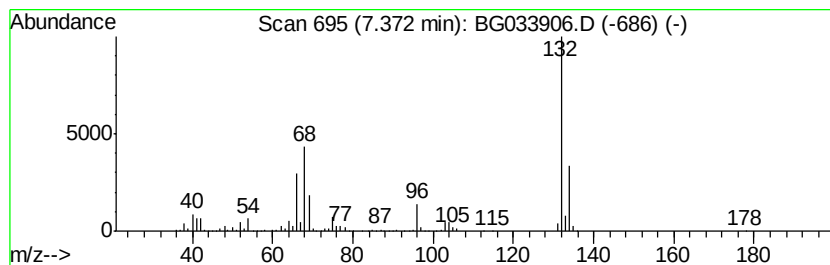
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Peak Number 2 unknown7.37 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.37	70.05 ng	1284750	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	35
2		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	16
3		Tranvlcyproline-propionyl	189	C12H15NO	1000123-86-3	10
4		Xenon	132	Xe	007440-63-3	9
5		Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	4.99	5.3	ng	96736	1	7.83	366814	20.0
unknown7.37	7.37	70.0	ng	1284750	1	7.83	366814	20.0