

Data Path : Z:\HPCHEM1\BNA G\DATA\BG042018\
 Data File : BG033905.D
 Acq On : 21 Apr 2018 4:27
 Operator : SJ/JU
 Sample : J2473-08
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 041818-SIDI-F-U-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	4.994	284	290	296	rBV	67848	103886	1.52%	0.432%
2	5.441	359	366	375	rBV	449367	688578	10.09%	2.862%
3	6.992	624	630	639	rVB	280170	475987	6.97%	1.978%
4	7.362	685	693	708	rBV	763292	1302524	19.08%	5.414%
5	7.832	765	773	779	rBV	237379	393924	5.77%	1.637%
6	8.149	819	827	836	rVB	988301	1669230	24.46%	6.938%
7	8.983	960	969	979	rBV	820576	1459586	21.38%	6.067%
8	10.617	1239	1247	1253	rBV	317818	566775	8.30%	2.356%
9	13.090	1660	1668	1679	rVB	2181599	3492293	51.16%	14.515%
10	14.465	1894	1902	1910	rBV2	495982	847047	12.41%	3.521%
11	15.952	2148	2155	2166	rBV2	1367075	1944441	28.49%	8.082%
12	17.215	2363	2370	2380	rVB2	760329	1124845	16.48%	4.675%
13	18.132	2522	2526	2540	rBV2	51782	94359	1.38%	0.392%
14	19.853	2813	2819	2827	rBV	5086482	6825716	100.00%	28.370%
15	21.169	3039	3043	3052	rBV5	34091	76374	1.12%	0.317%
16	21.469	3087	3094	3102	rBV	977008	1512762	22.16%	6.288%
17	24.536	3604	3616	3628	rVB	559509	1481312	21.70%	6.157%

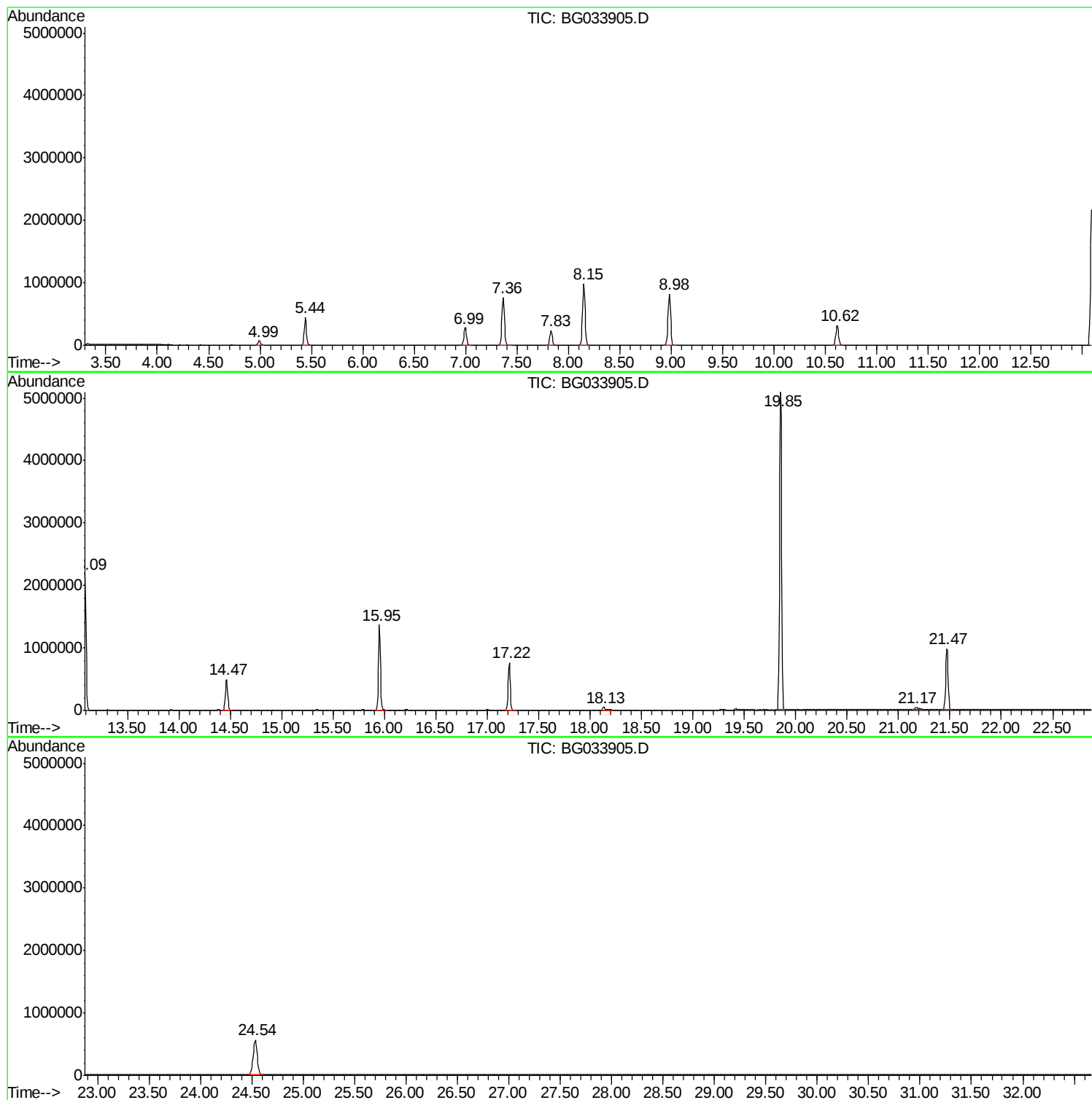
Sum of corrected areas: 24059639

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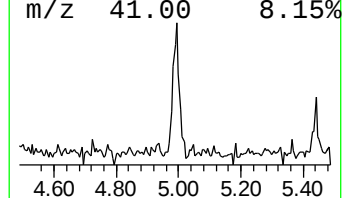
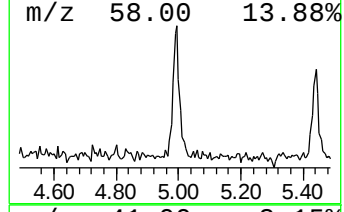
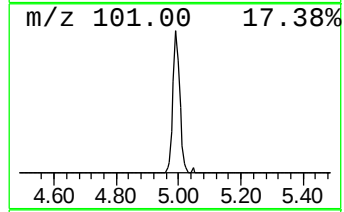
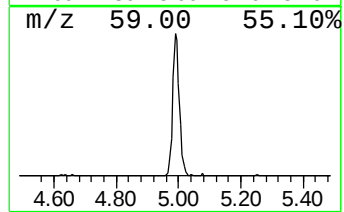
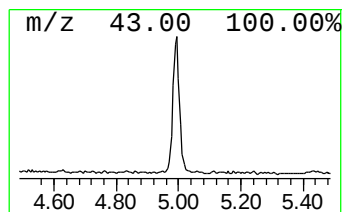
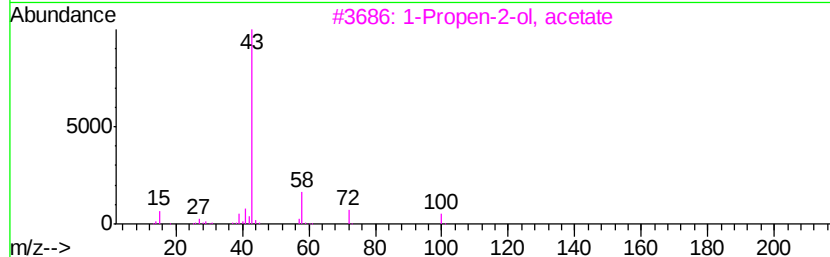
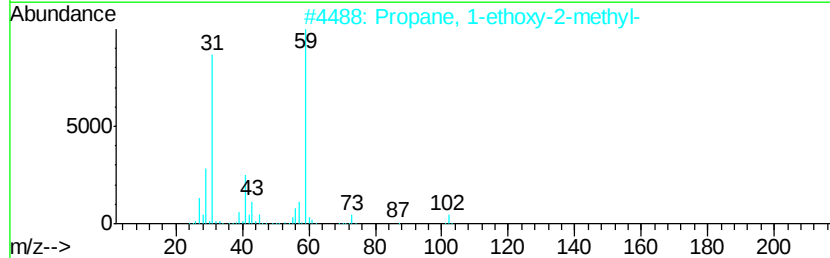
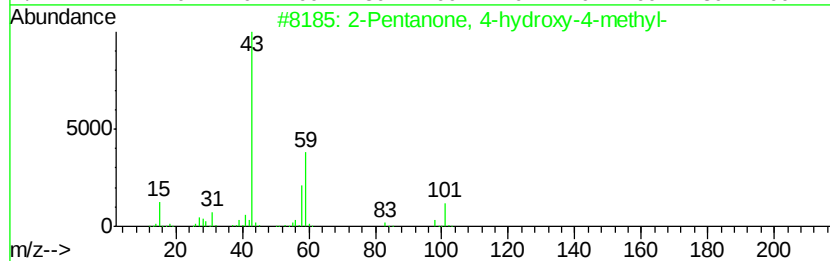
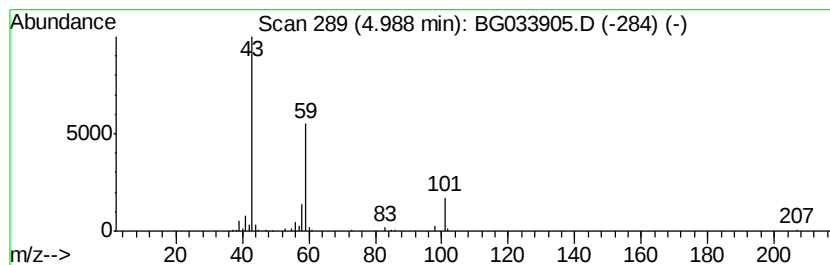
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.	
4.99	5.27 ng	103886	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, 1-ethoxy-2-methyl-	102	C6H14O	000627-02-1	32
3	1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	10
4	Silane, dimethyl-	60	C2H8Si	001111-74-6	9
5	Acetamide, N-(aminocarbonyl)-	102	C3H6N2O2	000591-07-1	9



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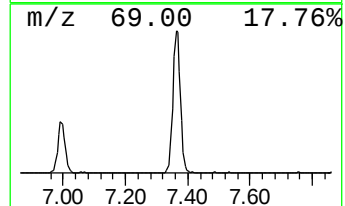
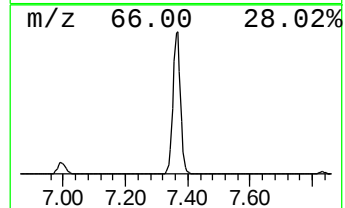
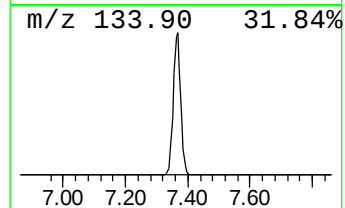
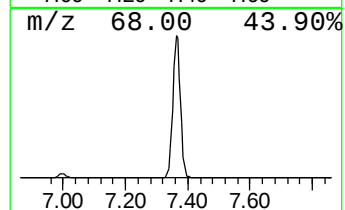
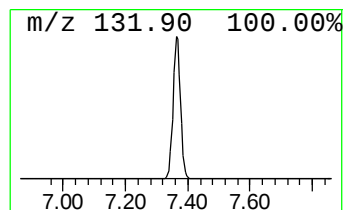
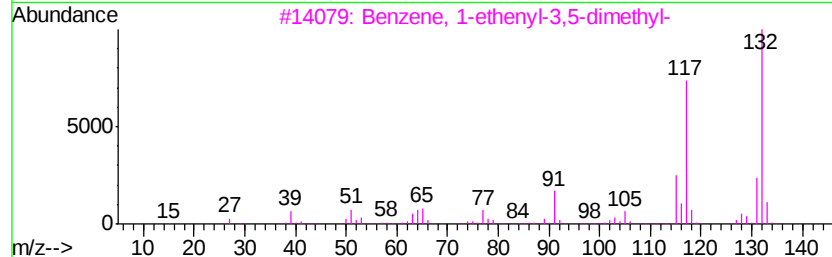
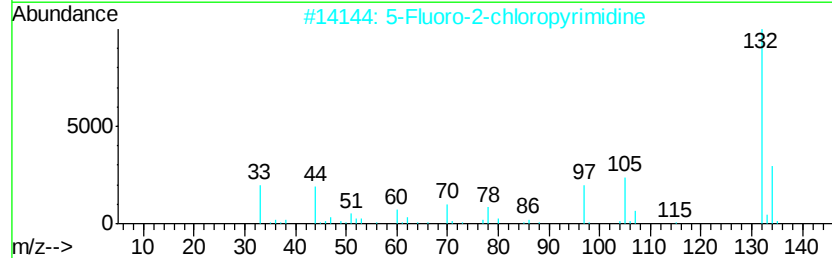
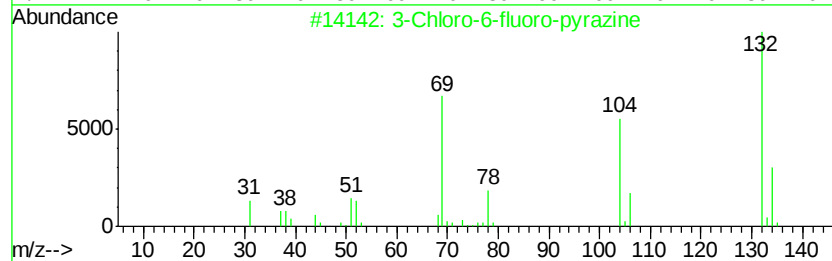
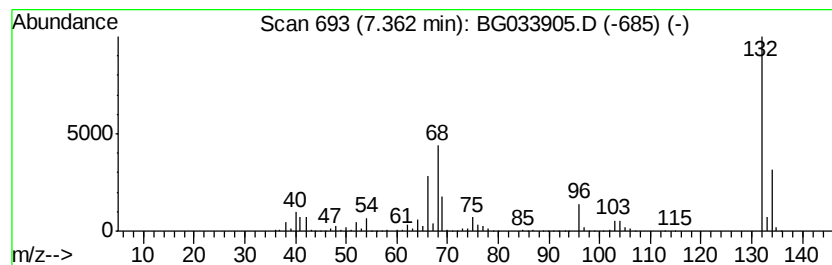
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Peak Number 2 unknown7.36 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.36	66.13 ng	1302520	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	16
2		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	12
3		Benzene, 1-ethenyl-3,5-dimethyl-	132	C10H12	005379-20-4	9
4		1-Methylpyrrolo[1,2-a]pyrazine	132	C8H8N2	064608-59-9	9
5		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	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	103886	1	7.83	393924	20.0
unknown7.36	7.36	66.1	ng	1302520	1	7.83	393924	20.0