

Data Path : Z:\HPCHEM1\BNA G\DATA\BG042518\
 Data File : BG033995.D
 Acq On : 25 Apr 2018 20:24
 Operator : SJ/JU
 Sample : PB108550BL
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB108550BL

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	4.986	283	289	298	rVB	61750	93844	1.40%	0.324%
2	5.438	358	366	380	rBV	1331734	2041301	30.50%	7.052%
3	7.001	622	632	643	rBV	1340787	2252692	33.65%	7.782%
4	7.365	685	694	707	rBV	1273319	2185901	32.66%	7.552%
5	7.830	766	773	782	rBV	215031	362858	5.42%	1.254%
6	8.147	819	827	837	rVB	908015	1560243	23.31%	5.390%
7	8.975	960	968	978	rBV	768806	1341947	20.05%	4.636%
8	10.609	1238	1246	1253	rBV	320086	563384	8.42%	1.946%
9	13.082	1657	1667	1677	rBV	2291339	3585290	53.56%	12.386%
10	14.457	1893	1901	1910	rBV2	553707	876034	13.09%	3.026%
11	15.949	2148	2155	2170	rBV	2072606	3094366	46.23%	10.690%
12	17.207	2361	2369	2377	rBV2	795522	1226737	18.33%	4.238%
13	19.845	2812	2818	2824	rBV	5080462	6693743	100.00%	23.124%
14	21.467	3087	3094	3101	rBV	997831	1582595	23.64%	5.467%
15	24.522	3603	3614	3626	rVB2	580853	1485622	22.19%	5.132%

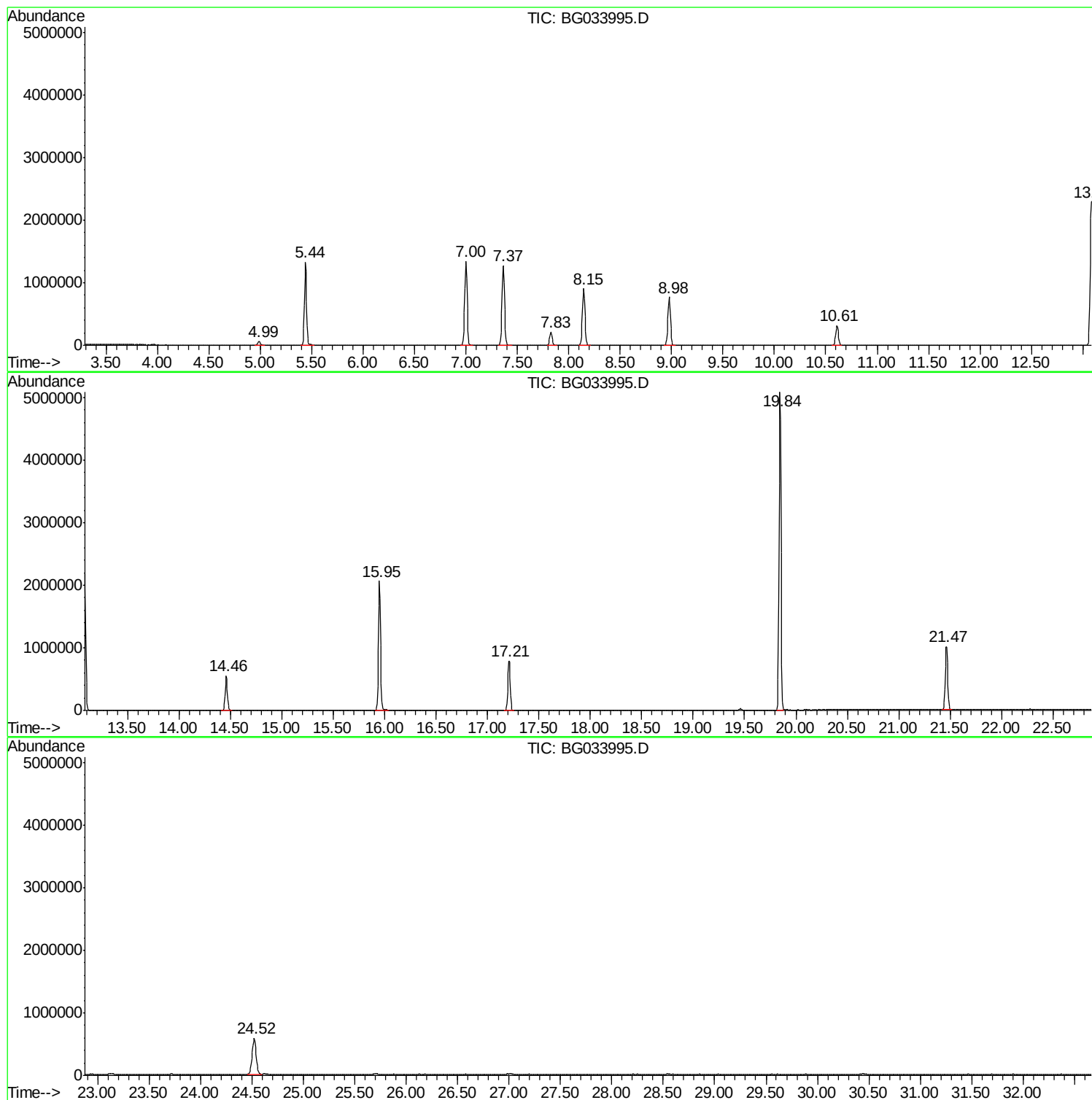
Sum of corrected areas: 28946557

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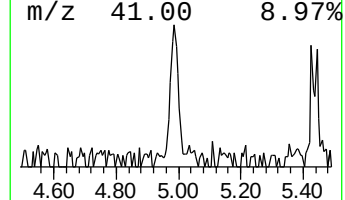
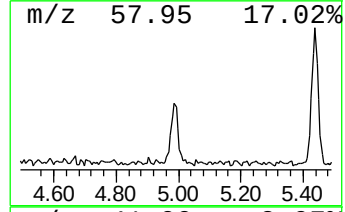
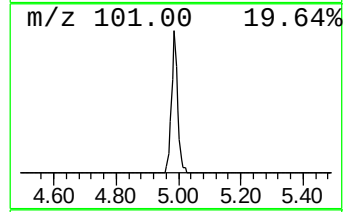
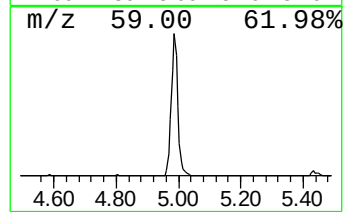
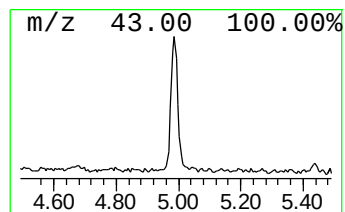
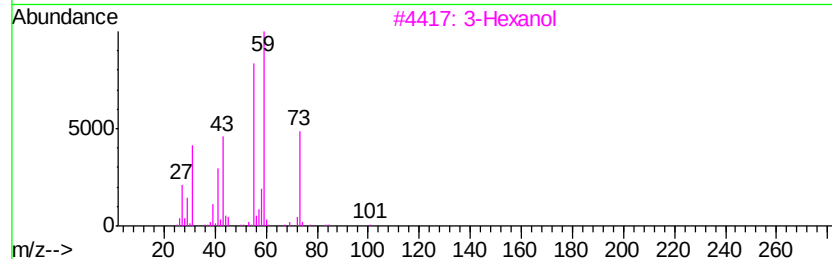
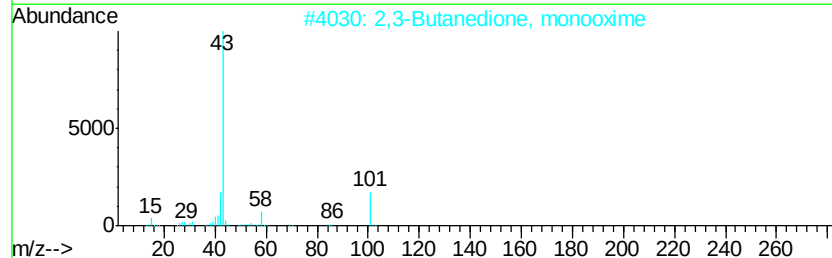
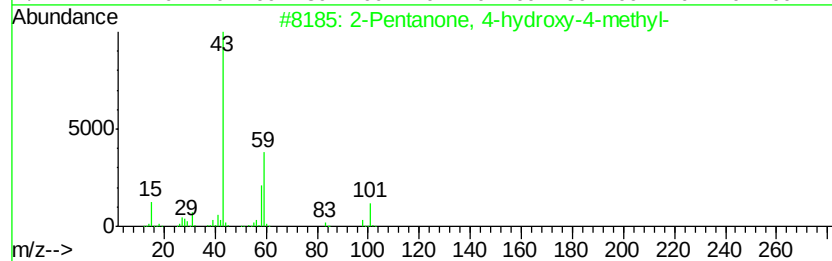
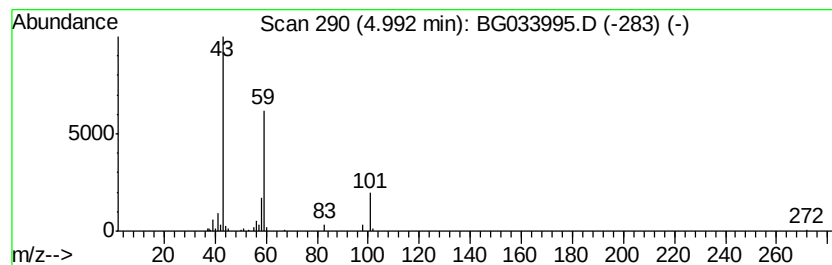
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.99	5.17 ng	93844	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	2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	25	
3	3-Hexanol	102	C6H14O	000623-37-0	25	
4	Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	25	
5	Propanoic acid, 3-nitro-, methyl...	133	C4H7NO4	020497-95-4	17	



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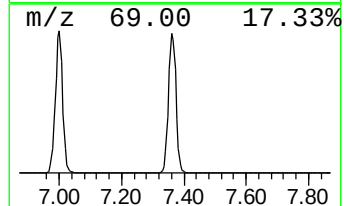
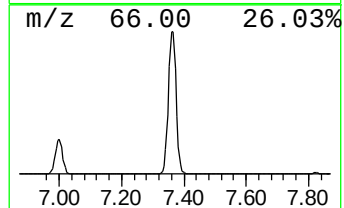
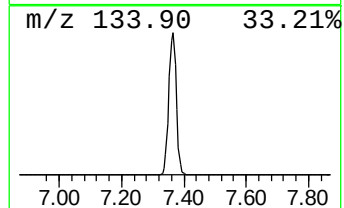
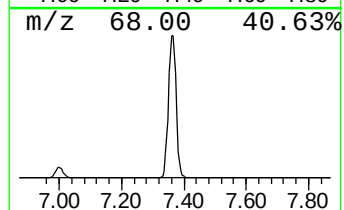
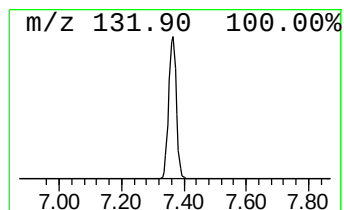
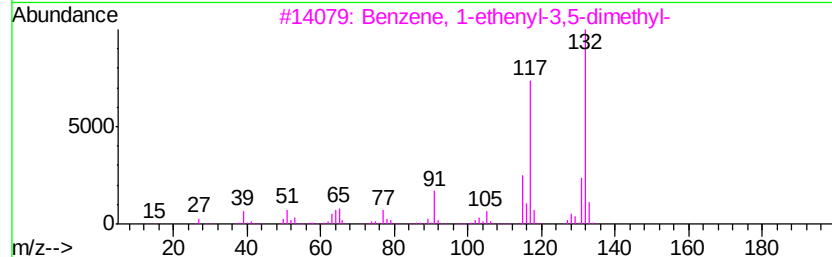
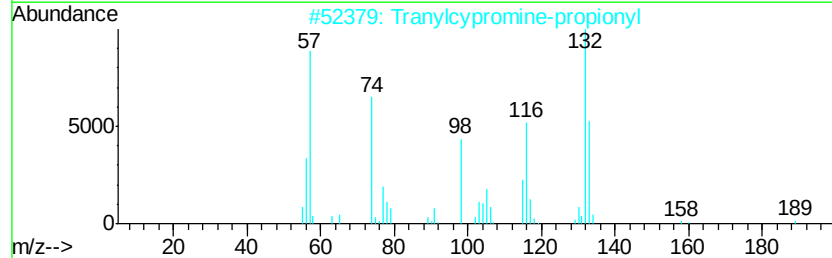
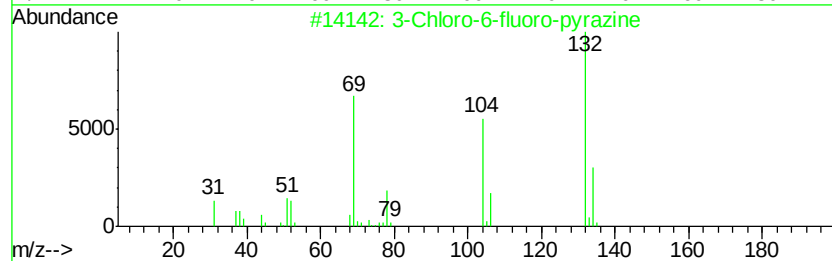
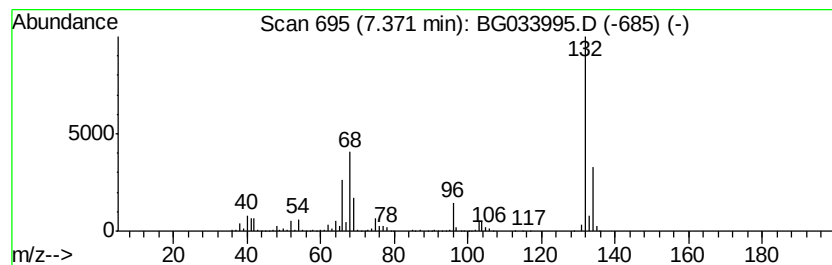
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Peak Number 2 unknown7.37 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.37	120.48 ng	2185900	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		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
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		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	4.99	5.2	ng	93844	1	7.83	362858	20.0
unknown7.37	7.37	120.5	ng	2185900	1	7.83	362858	20.0