

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG091417\
 Data File : BG028730.D
 Acq On : 14 Sep 2017 13:24
 Operator : SJ/JU
 Sample : PB102225BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :

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-BG091217.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	5.435	308	314	322	rBV2	37141	62060	2.00%	0.522%
2	5.899	386	393	406	rBV	496145	831181	26.77%	6.996%
3	7.480	652	662	674	rBV	453055	844100	27.18%	7.105%
4	7.862	719	727	736	rBV	459409	819750	26.40%	6.900%
5	8.326	799	806	814	rBV	83056	145642	4.69%	1.226%
6	8.655	854	862	873	rVB	339275	611314	19.69%	5.146%
7	9.495	997	1005	1014	rBV	255769	497160	16.01%	4.185%
8	11.146	1278	1286	1295	rBV	97569	196658	6.33%	1.655%
9	13.561	1690	1697	1710	rBV	802658	1312727	42.28%	11.050%
10	14.942	1925	1932	1940	rBV	179873	313013	10.08%	2.635%
11	16.428	2179	2185	2196	rBV	785837	1345602	43.34%	11.327%
12	17.691	2390	2400	2413	rBV	317265	513533	16.54%	4.323%
13	20.276	2833	2840	2849	rBV	2297788	3105087	100.00%	26.137%
14	22.021	3130	3137	3150	rBV	368845	665963	21.45%	5.606%
15	25.523	3722	3733	3752	rBV	171630	616295	19.85%	5.188%

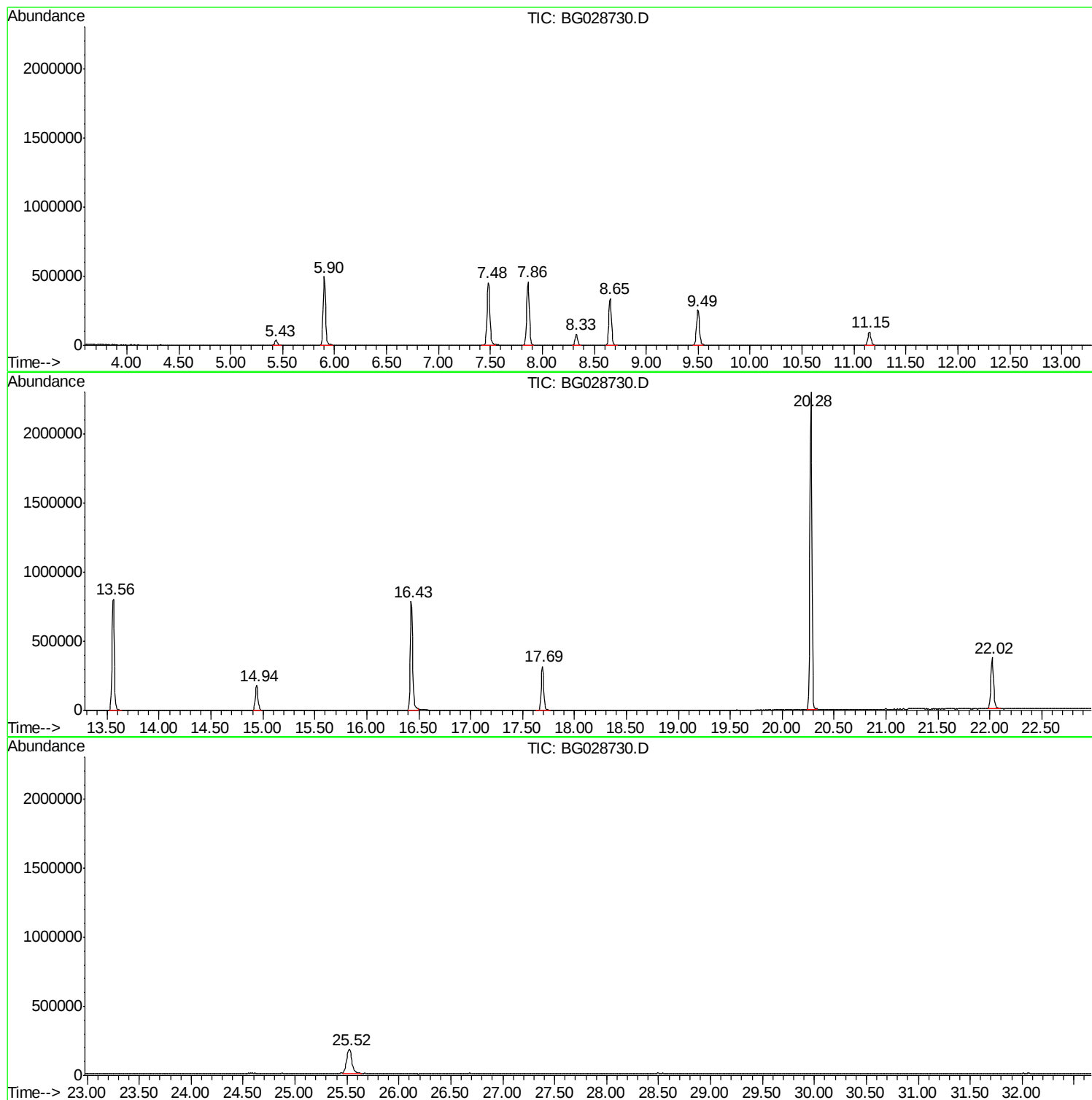
Sum of corrected areas: 11880085

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Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG091217.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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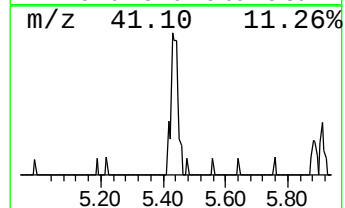
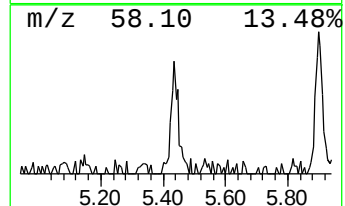
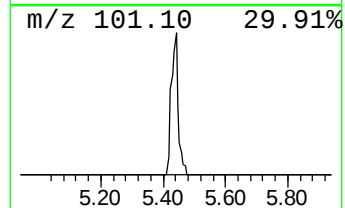
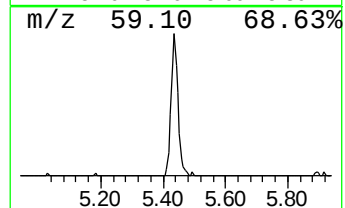
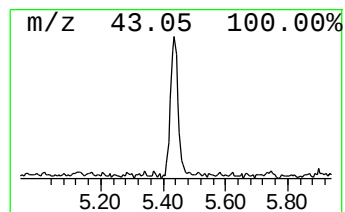
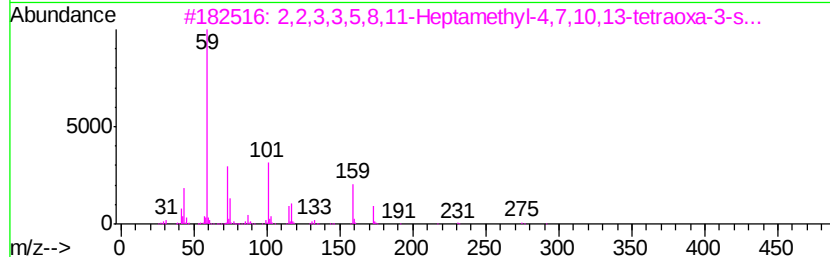
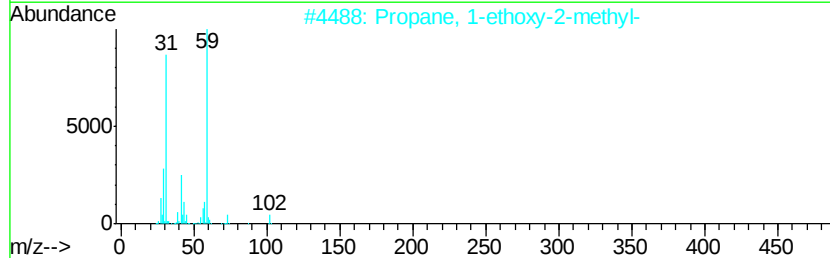
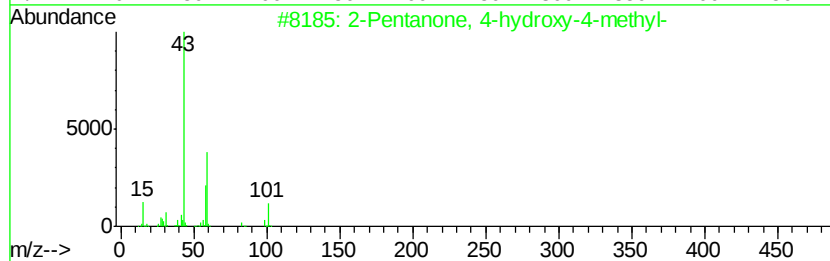
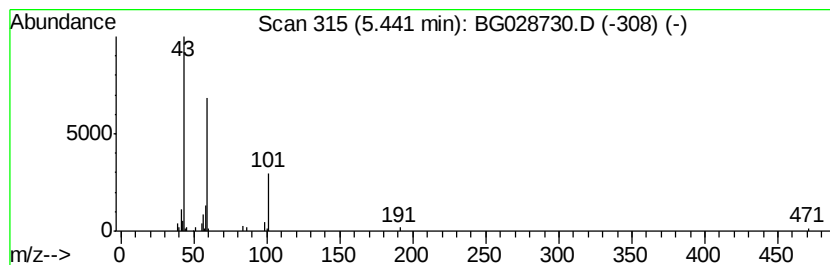
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.44	8.52 ng	62060	1,4-Dichlorobenzene-d4	8.33

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	43	
3	2,2,3,3,5,8,11-Heptamethyl-4,7,1...	348	C18H40O4Si	1000367-07-2	36	
4	n-Propyl t-butyl ether	116	C7H16O	1000334-81-9	36	
5	3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	28	



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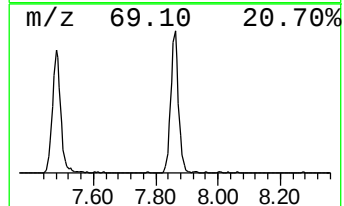
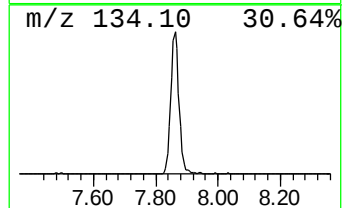
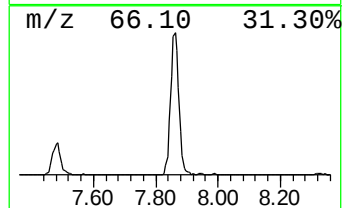
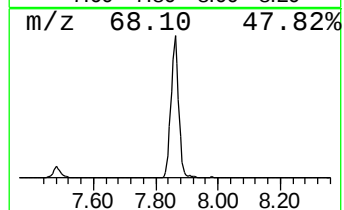
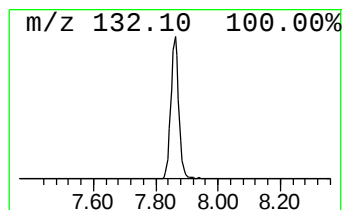
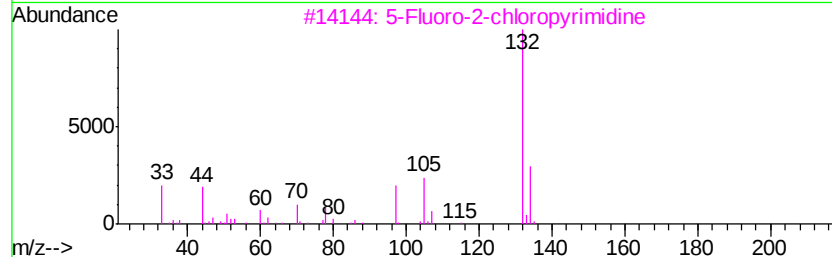
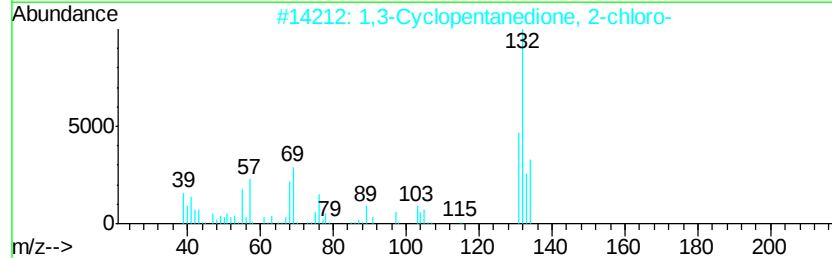
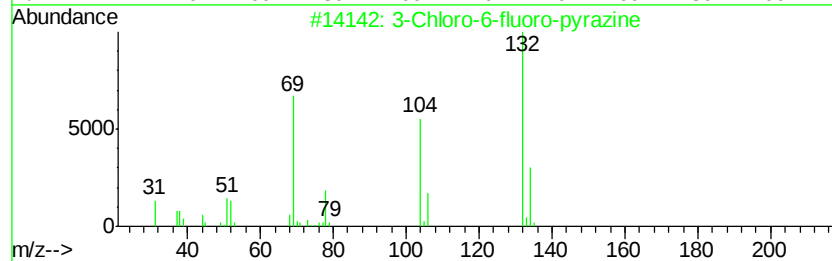
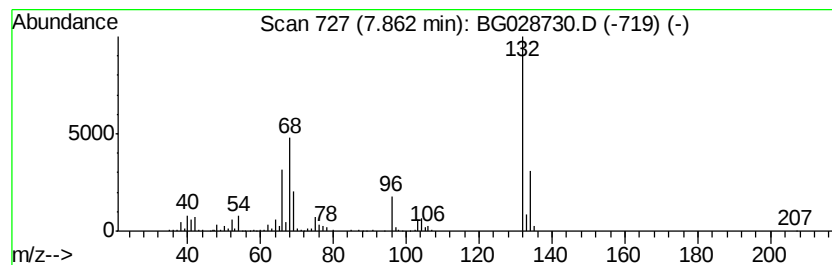
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Peak Number 2 unknown7.86 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.86	112.57 ng	819750	1,4-Dichlorobenzene-d4	8.33

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		1,3-Cyclopentanedione, 2-chloro-	132	C5H5ClO2	014203-19-1	13
3		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	12
4		5-Aminoindole	132	C8H8N2	005192-03-0	12
5		Cyclopropanamine, 1-phenyl-	133	C9H11N	1000351-26-2	10



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.44	8.5	ng	62060	1	8.33	145642	20.0
unknown7.86	7.86	112.6	ng	819750	1	8.33	145642	20.0