

Data Path : Z:\HPCHEM1\BNA G\DATA\BG092517\
 Data File : BG028936.D
 Acq On : 26 Sep 2017 9:51
 Operator : SJ/JU
 Sample : I5385-02
 Misc :
 ALS Vial : 27 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 092117-SIDI-F-D-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-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	5.380	300	305	313	rBV2	31581	49428	1.29%	0.432%
2	5.844	377	384	393	rBV	163827	284171	7.42%	2.484%
3	7.436	647	655	670	rBV	89173	186497	4.87%	1.630%
4	7.812	711	719	728	rBV	285701	521231	13.61%	4.556%
5	8.276	792	798	806	rBV2	84681	146010	3.81%	1.276%
6	8.605	844	854	865	rBV	382640	720932	18.83%	6.302%
7	9.446	989	997	1011	rBV	313230	621706	16.24%	5.435%
8	11.097	1270	1278	1286	rBV	104897	198166	5.18%	1.732%
9	13.512	1679	1689	1698	rBV	997434	1618638	42.28%	14.150%
10	14.892	1918	1924	1936	rVB3	176045	316782	8.27%	2.769%
11	16.385	2170	2178	2192	rBV	673859	1092437	28.53%	9.550%
12	17.648	2387	2393	2399	rBV2	318218	529051	13.82%	4.625%
13	20.233	2827	2833	2845	rBV	2852200	3828614	100.00%	33.468%
14	21.960	3120	3127	3136	rBV	374637	704513	18.40%	6.159%
15	25.421	3706	3716	3727	rBV2	200940	621338	16.23%	5.432%

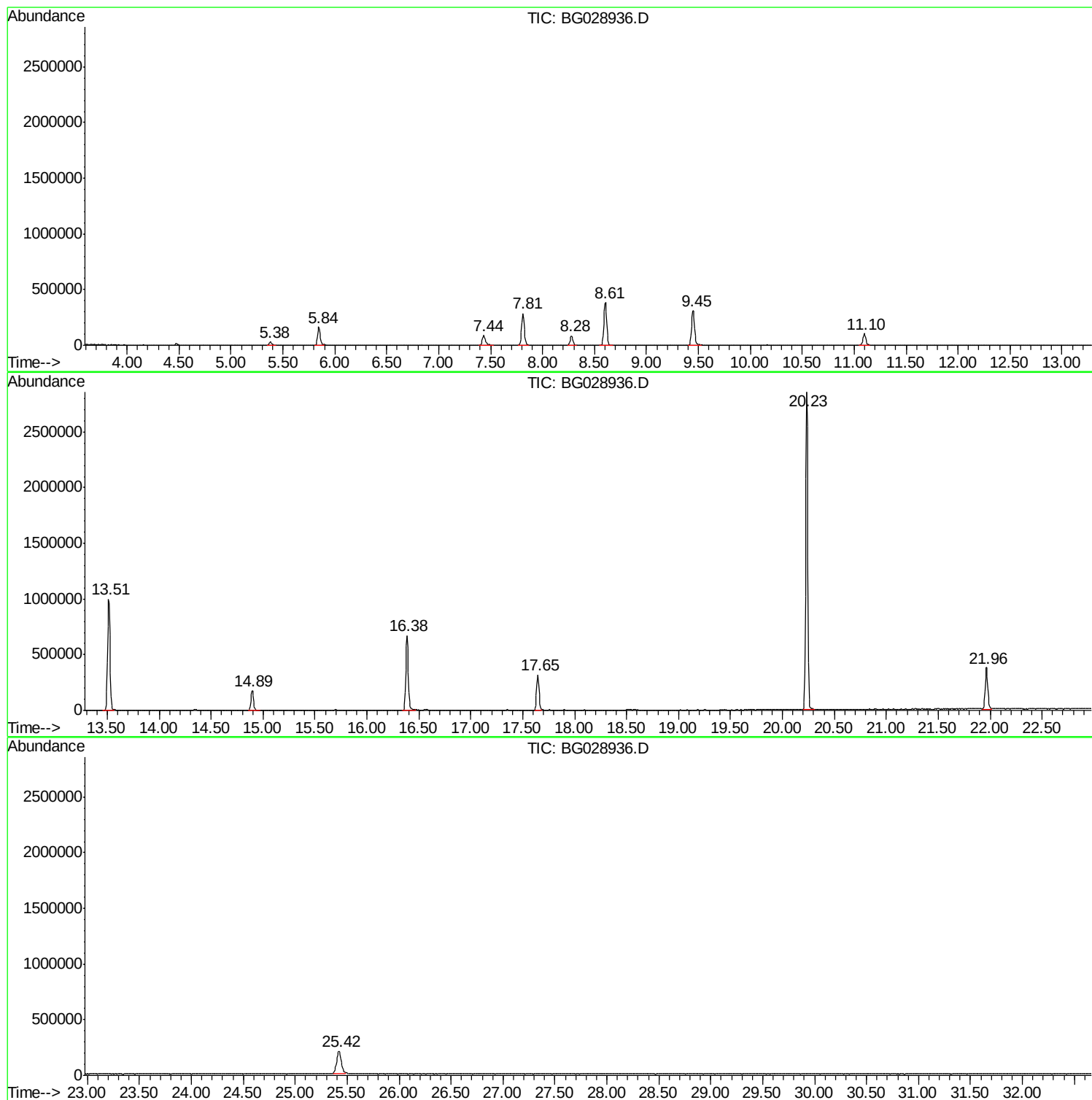
Sum of corrected areas: 11439514

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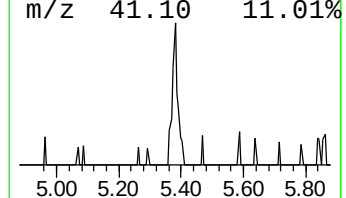
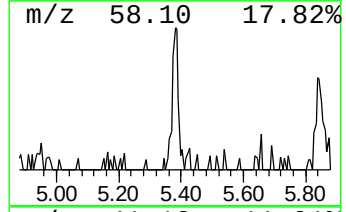
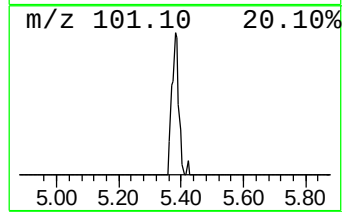
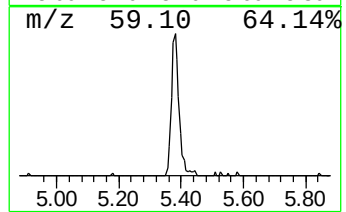
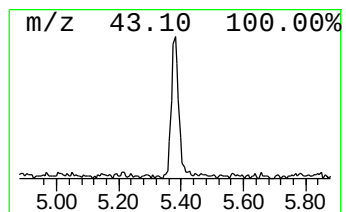
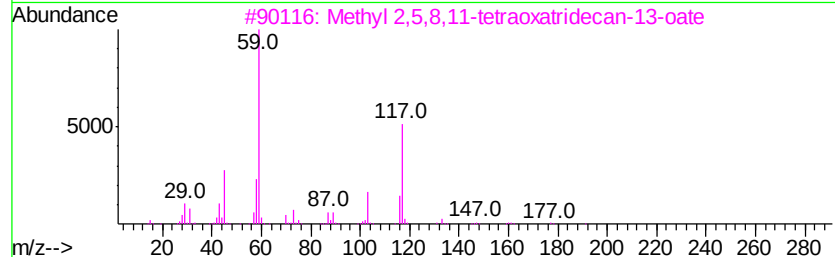
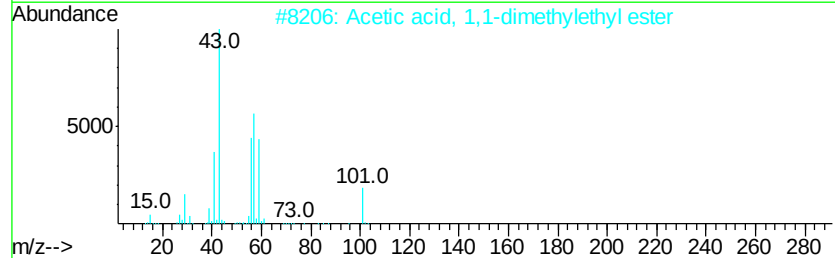
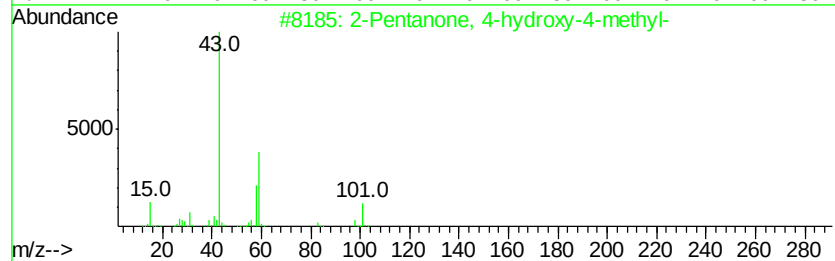
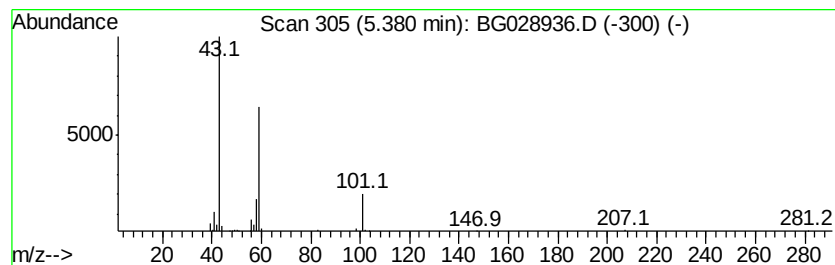
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.38	6.77 ng	49428	1,4-Dichlorobenzene-d4	8.28

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	59
2			Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	38
3			Methyl 2,5,8,11-tetraoxatridecan...	236	C10H20O6	1000366-78-2	23
4			3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	17
5			Propanoic acid, 3-nitro-, methyl...	133	C4H7NO4	020497-95-4	12



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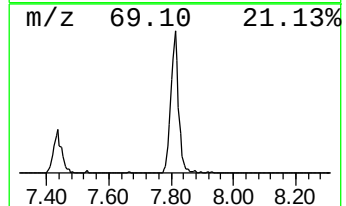
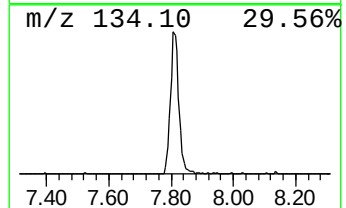
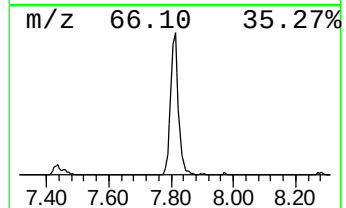
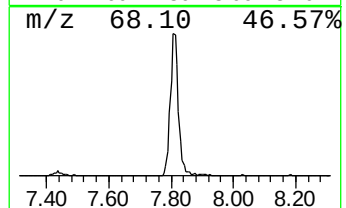
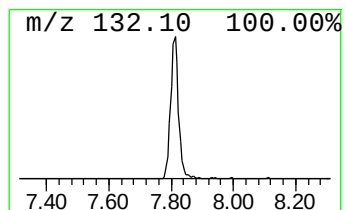
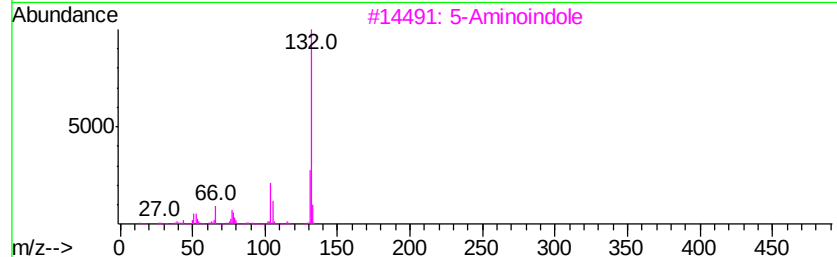
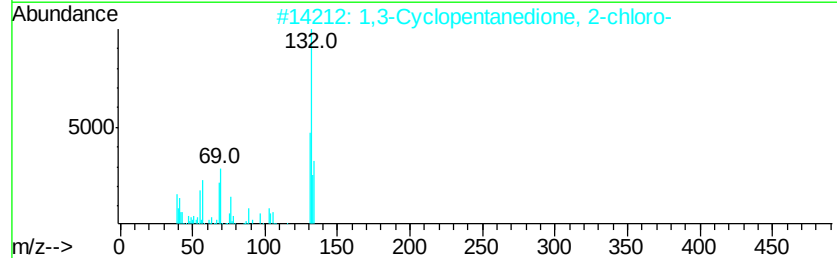
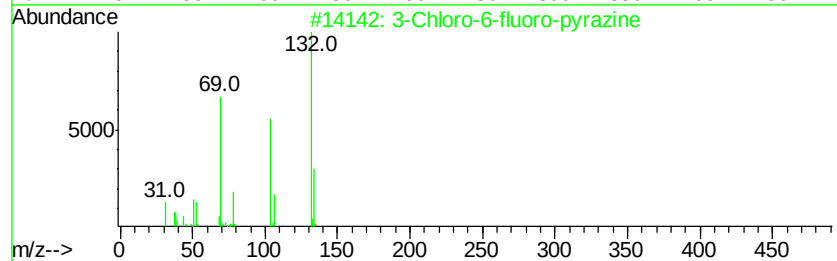
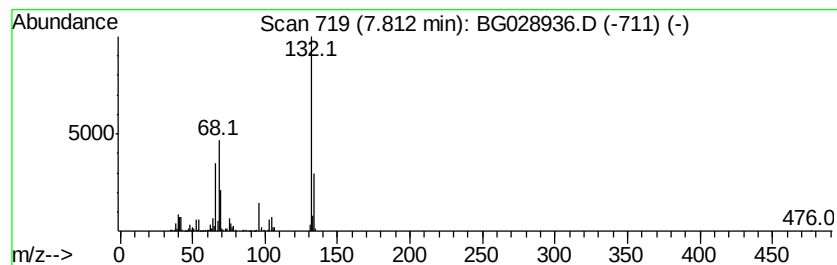
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Peak Number 2 unknown7.81 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.81	71.40 ng	521231	1,4-Dichlorobenzene-d4	8.28

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	16
2		1,3-Cyclopentanedione, 2-chloro-	132	C5H5ClO2	014203-19-1	13
3		5-Aminoindole	132	C8H8N2	005192-03-0	12
4		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	12
5		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.38	6.8	ng	49428	1	8.28	146010	20.0
unknown7.81	7.81	71.4	ng	521231	1	8.28	146010	20.0