

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG022417\
 Data File : BG025960.D
 Acq On : 25 Feb 2017 9:38
 Operator : SJ/MA
 Sample : I1969-01
 Misc :
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 022217-PRRT-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-BG021717.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	5.176	306	313	321	rBV	139327	213619	2.49%	0.660%
2	5.646	387	393	402	rBV	442177	688369	8.02%	2.126%
3	7.232	656	663	675	rBV	271592	469040	5.47%	1.449%
4	7.608	718	727	736	rBV	811336	1398357	16.30%	4.319%
5	8.078	800	807	815	rBV	270398	466249	5.43%	1.440%
6	8.407	854	863	873	rBV	1037618	1813348	21.14%	5.601%
7	9.236	995	1004	1013	rBV	936701	1662176	19.38%	5.134%
8	10.881	1277	1284	1292	rVB	461815	802999	9.36%	2.480%
9	13.313	1689	1698	1707	rBV	3065426	4775939	55.67%	14.752%
10	14.124	1829	1836	1842	rBV	249847	361844	4.22%	1.118%
11	14.682	1923	1931	1942	rVB2	827936	1343951	15.67%	4.151%
12	16.157	2175	2182	2192	rBV	1958517	2989373	34.85%	9.233%
13	17.414	2389	2396	2404	rBV	1276873	1965977	22.92%	6.072%
14	18.290	2540	2545	2553	rBV2	96814	160934	1.88%	0.497%
15	19.553	2755	2760	2769	rBV2	61923	105004	1.22%	0.324%
16	20.011	2831	2838	2844	rBV	6343390	8578820	100.00%	26.498%
17	21.298	3053	3057	3067	rBV	125066	191728	2.23%	0.592%
18	21.662	3111	3119	3126	rBV	1296049	2195828	25.60%	6.782%
19	24.865	3653	3664	3678	rBV	798325	2192308	25.55%	6.771%

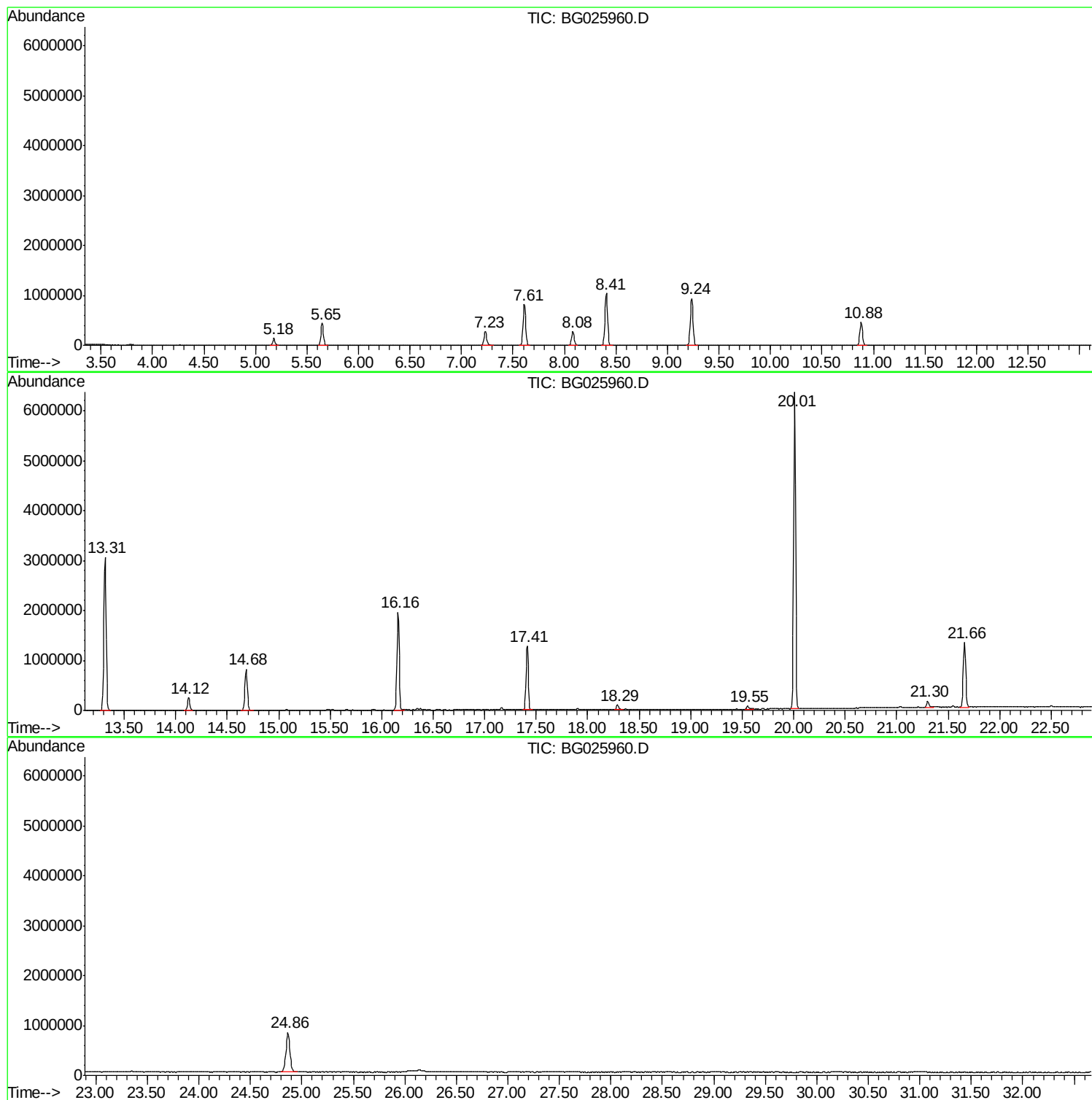
Sum of corrected areas: 32375863

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Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG021717.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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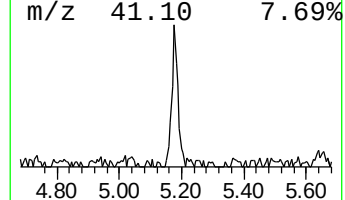
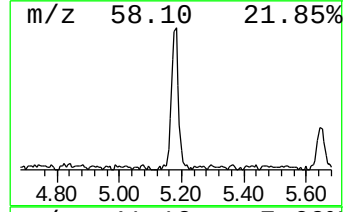
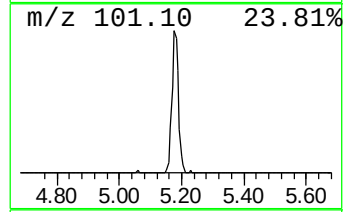
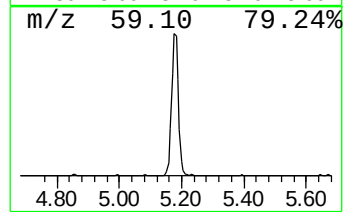
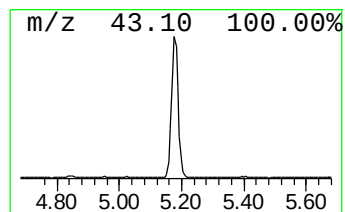
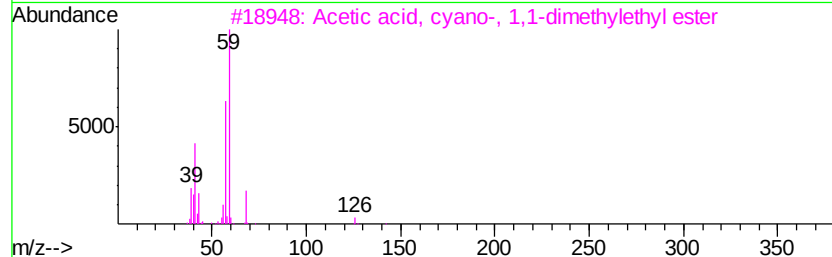
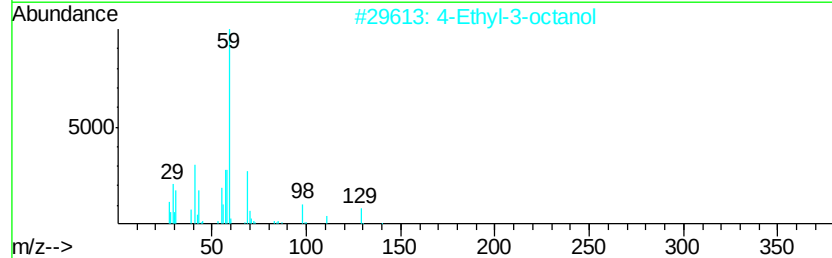
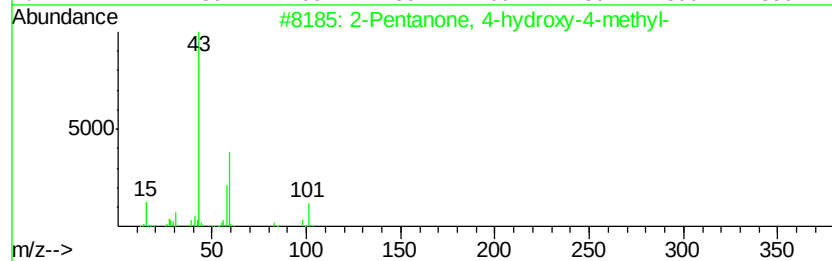
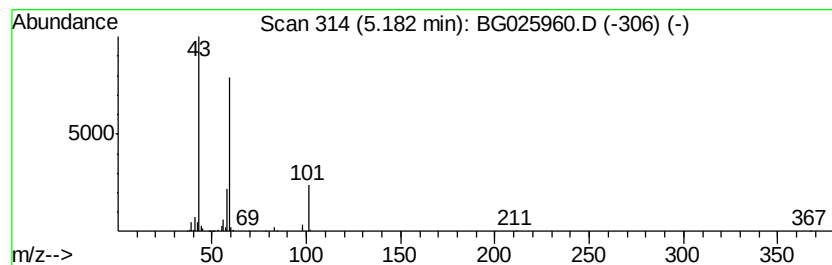
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.18	9.16 ng	213619	1,4-Dichlorobenzene-d4	8.08

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	38
2			4-Ethyl-3-octanol	158	C10H22O	019781-26-1	33
3			Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	28
4			Propane, 1-ethoxy-2-methyl-	102	C6H14O	000627-02-1	25
5			1,8-Nonanediol, 8-methyl-	174	C10H22O2	054725-73-4	25



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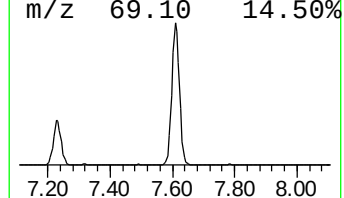
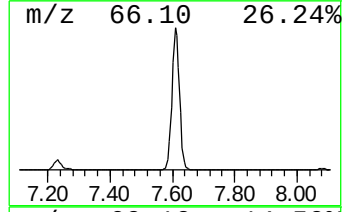
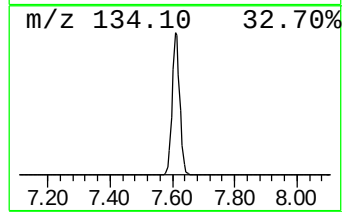
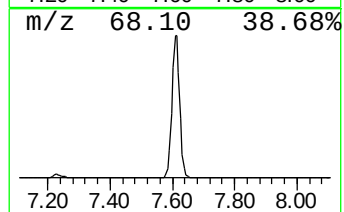
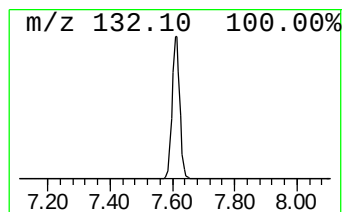
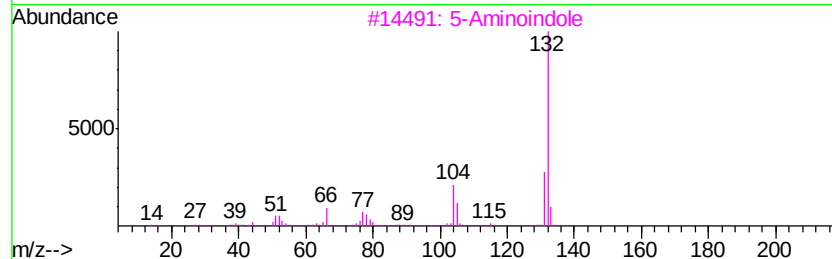
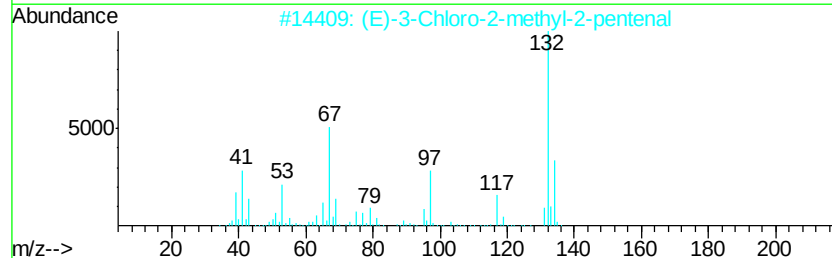
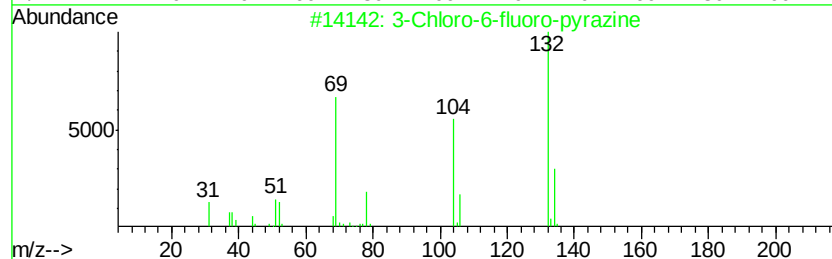
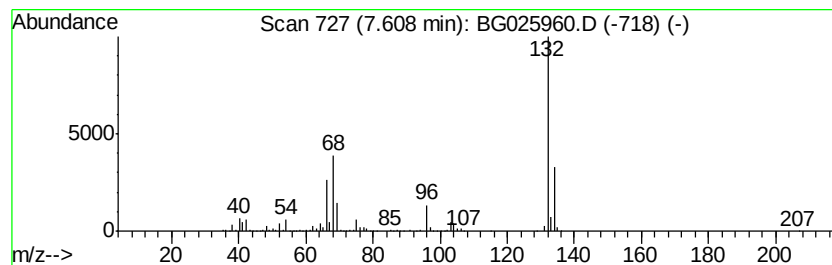
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Peak Number 2 unknown7.61 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.61	59.98 ng	1398360	1,4-Dichlorobenzene-d4	8.08

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	16
3		5-Aminoindole	132	C8H8N2	005192-03-0	12
4		Tranlylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
5		1,3,5-Trifluorobenzene	132	C6H3F3	000372-38-3	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.18	9.2	ng	213619	1	8.08	466249	20.0
unknown7.61	7.61	60.0	ng	1398360	1	8.08	466249	20.0