

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG110216\
Data File : BG024636.D
Acq On : 2 Nov 2016 21:54
Operator : UM/SJ
Sample : H5491-08
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
GW-69-7.6-12.0

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-BG102516.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	5.331	416	424	433	rVB	366137	606160	7.07%	1.559%
2	5.801	496	504	512	rBV	854830	1481822	17.28%	3.812%
3	7.400	770	776	787	rVB	564146	1007839	11.75%	2.593%
4	7.793	833	843	854	rVB	1528139	2695417	31.43%	6.934%
5	8.269	917	924	935	rVB	430831	797513	9.30%	2.052%
6	8.592	968	979	989	rVB	1403887	2535954	29.57%	6.524%
7	9.438	1114	1123	1135	rBV	1190397	2209010	25.76%	5.683%
8	11.095	1395	1405	1412	rBV	568708	1057345	12.33%	2.720%
9	12.000	1552	1559	1570	rVB5	40633	103960	1.21%	0.267%
10	13.504	1806	1815	1824	rBV	3010871	4872131	56.81%	12.534%
11	14.879	2043	2049	2058	rBV	914093	1473259	17.18%	3.790%
12	16.366	2294	2302	2313	rBV	2523871	4213787	49.13%	10.841%
13	17.623	2509	2516	2526	rBV2	1187010	1889867	22.03%	4.862%
14	20.202	2949	2955	2961	rBV	6442997	8576874	100.00%	22.065%
15	21.912	3239	3246	3254	rVB	1521086	2721454	31.73%	7.001%
16	25.337	3818	3829	3842	rBV2	809561	2628134	30.64%	6.761%

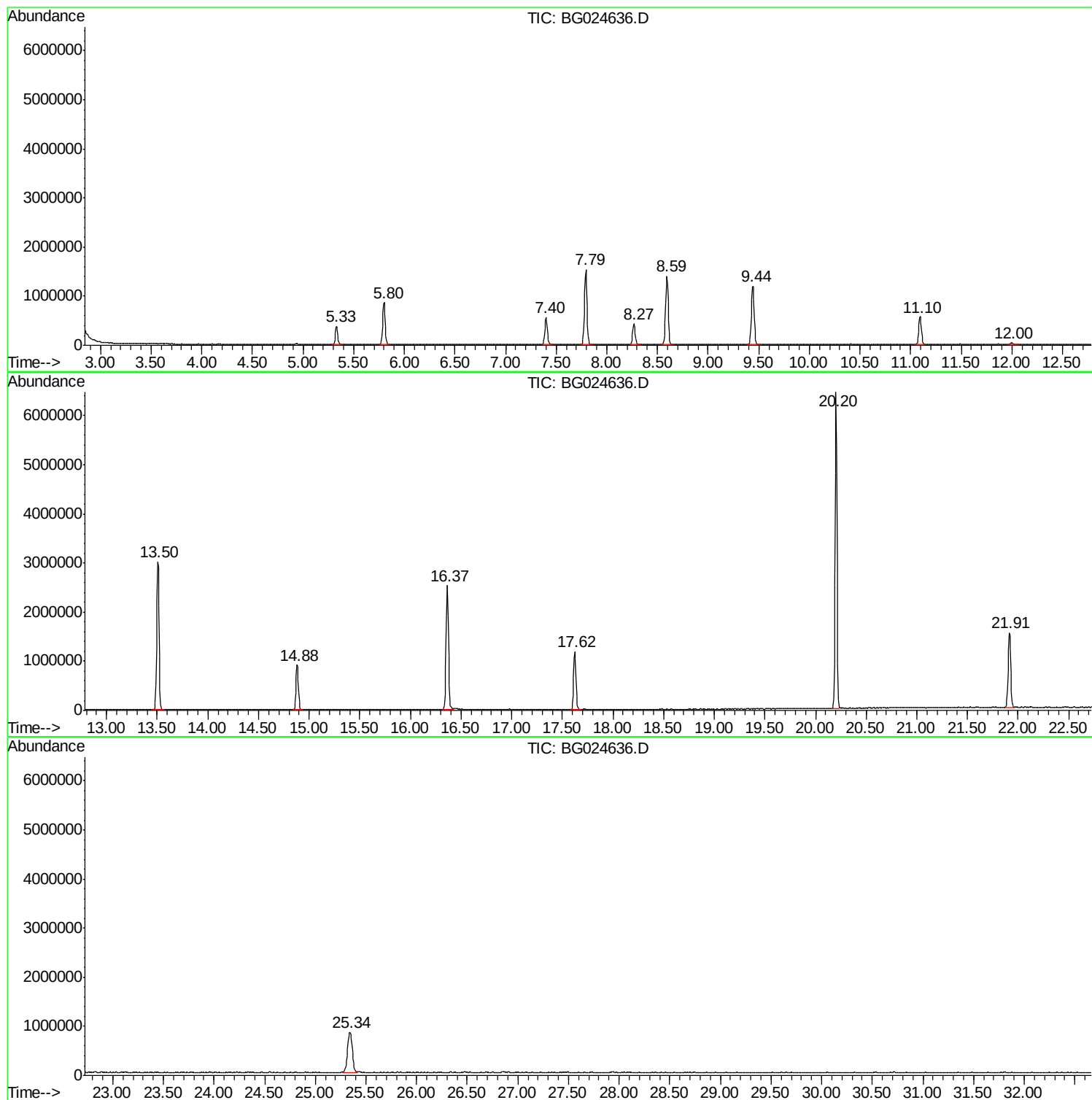
Sum of corrected areas: 38870526

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Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG102516.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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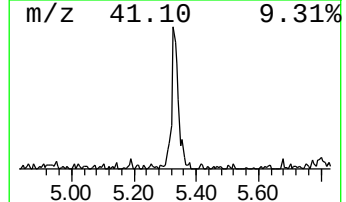
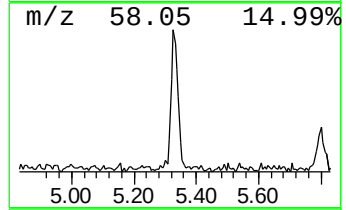
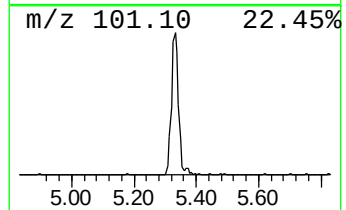
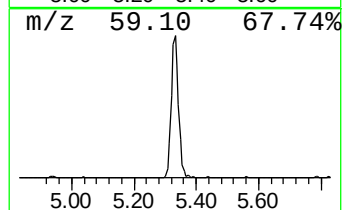
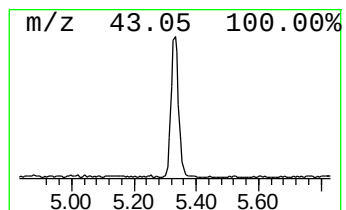
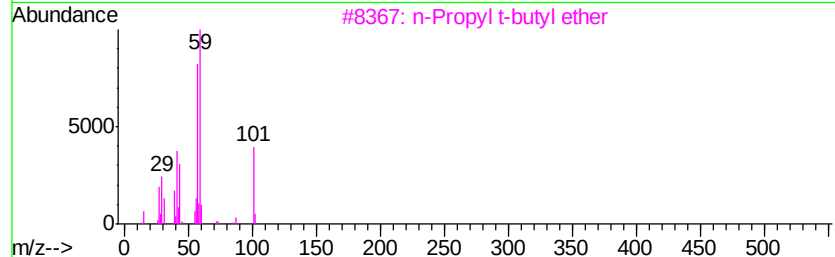
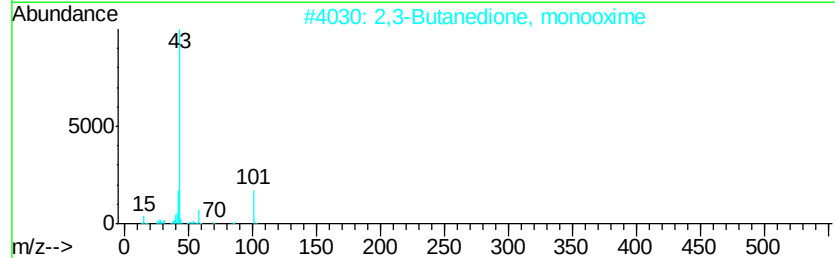
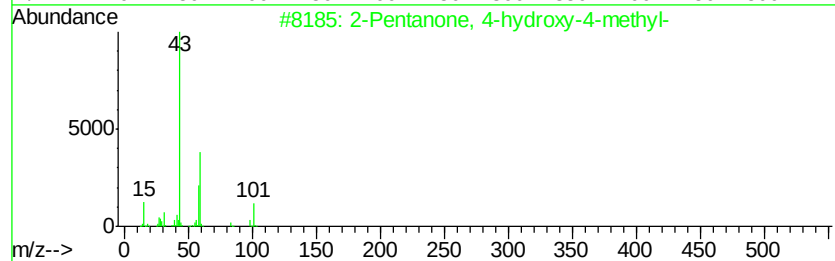
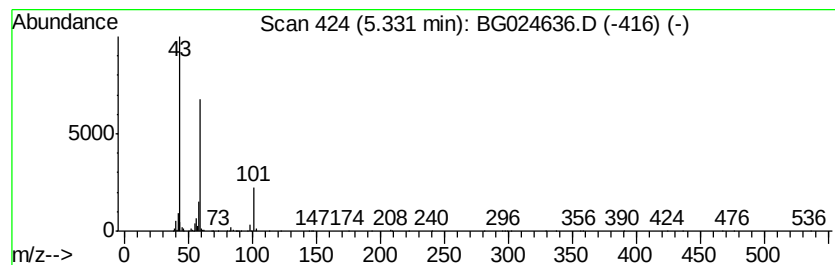
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.33	15.20 ng	606160	1,4-Dichlorobenzene-d4	8.27

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	45
2			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	32
3			n-Propyl t-butyl ether	116	C7H16O	1000334-81-9	28
4			Guanidine	59	CH5N3	000113-00-8	9
5			Butane, 1-ethoxy-	102	C6H14O	000628-81-9	9



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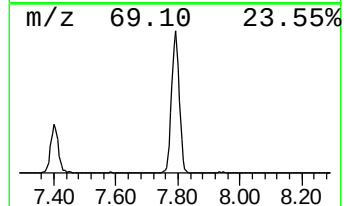
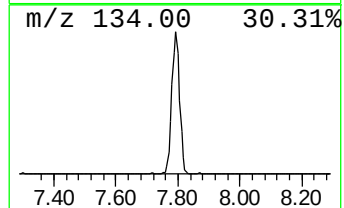
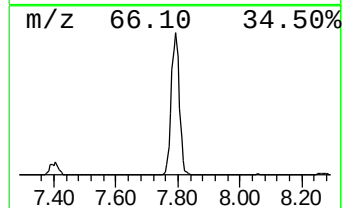
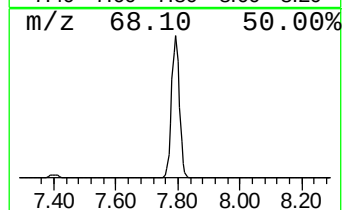
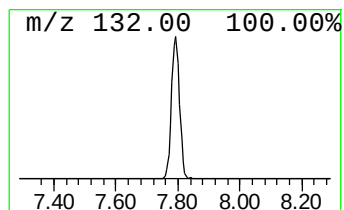
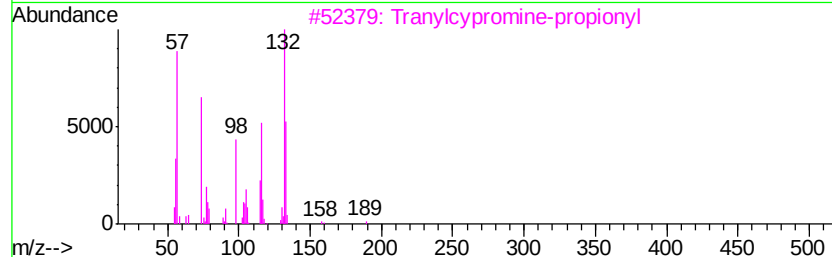
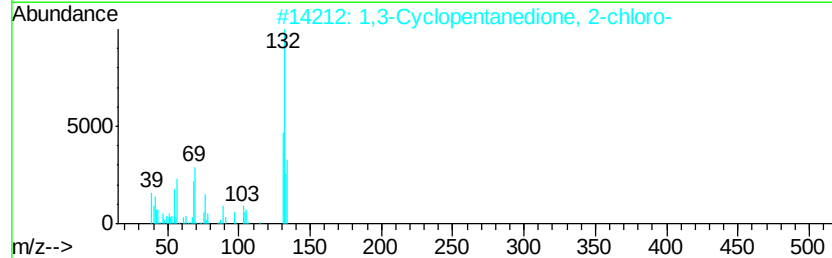
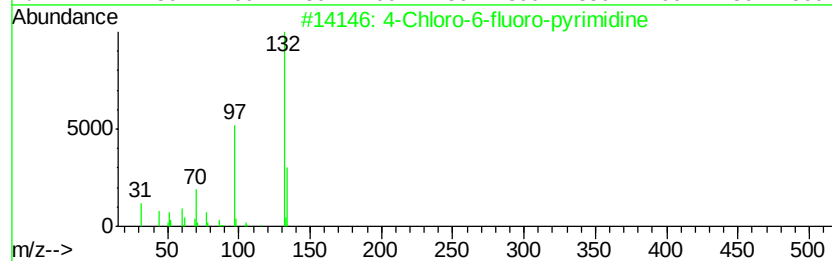
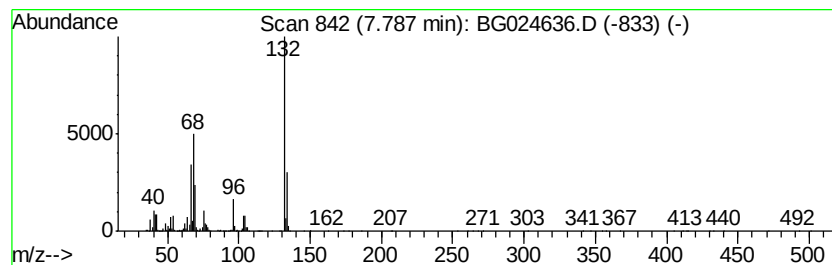
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Peak Number 2 unknown7.79 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.79	67.60 ng	2695420	1,4-Dichlorobenzene-d4	8.27

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	4-Chloro-6-fluoro-pyrimidine	132	C4H2ClFN2	051422-01-6	27
2		1,3-Cyclopentanedione, 2-chloro-	132	C5H5ClO2	014203-19-1	23
3		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	22
4		1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	22
5		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	16



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.33	15.2	ng	606160	1	8.27	797513	20.0
unknown7.79	7.79	67.6	ng	2695420	1	8.27	797513	20.0