

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF060515\
Data File : BF079751.D
Acq On : 6 Jun 2015 3:04
Operator : TP/IZ
Sample : G2464-13 10X
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SB-17AS

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_F\METHODS\8270-BF060515.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	1.324	10	12	14	rVB2	16306701	21425990	100.00%	76.511%
2	2.410	104	107	108	rBV	480385	662304	3.09%	2.365%
3	2.433	108	109	144	rVB	1435877	2779878	12.97%	9.927%
4	7.485	549	551	553	rBV	243625	282266	1.32%	1.008%
5	8.776	662	664	667	rBV	292100	391234	1.83%	1.397%
6	10.559	818	820	822	rVB	595607	495090	2.31%	1.768%
7	12.068	950	952	953	rBV	585120	538193	2.51%	1.922%
8	14.925	1199	1202	1203	rBV	552415	739953	3.45%	2.642%
9	16.811	1364	1367	1375	rVB	389472	688827	3.21%	2.460%

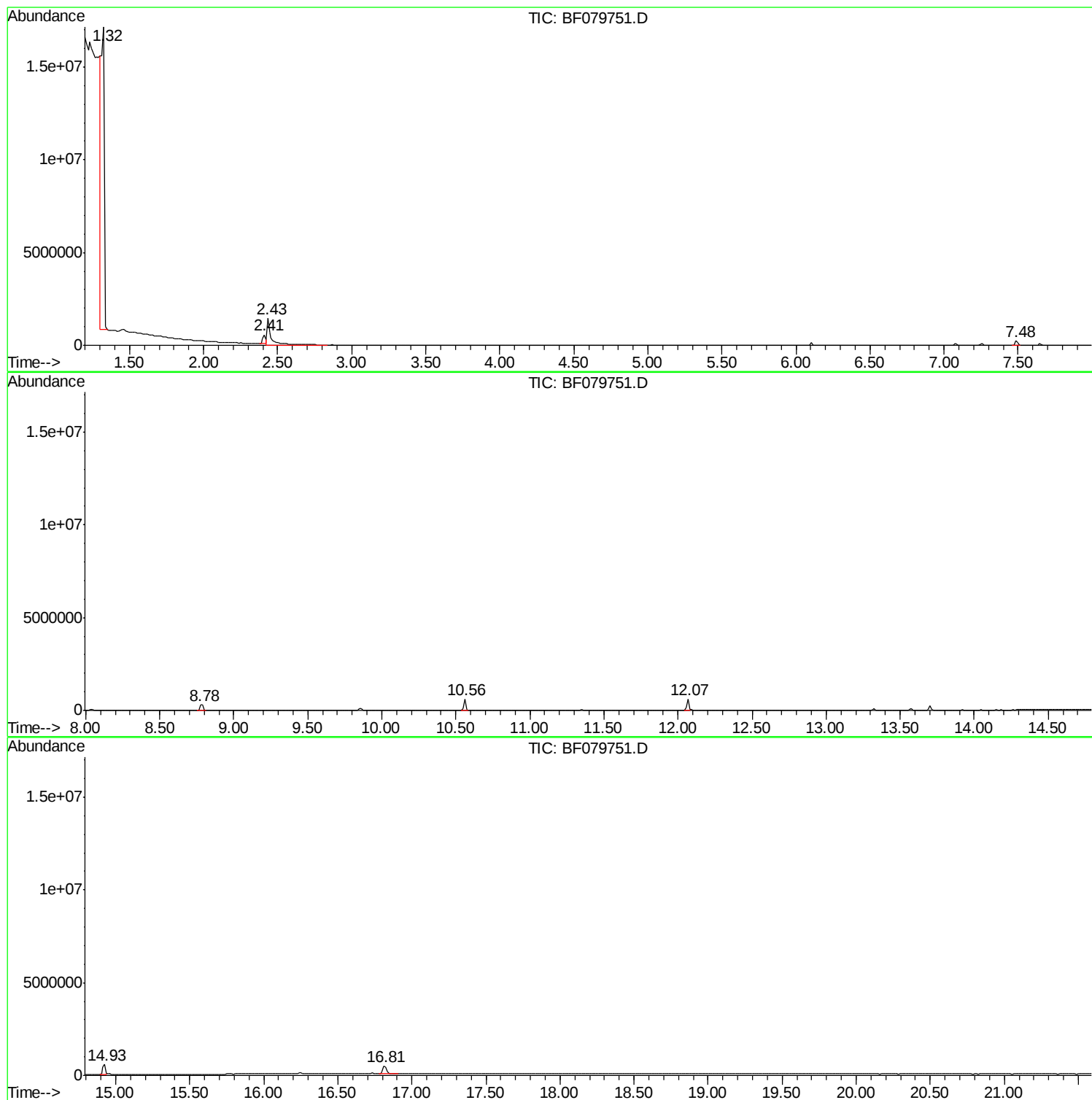
Sum of corrected areas: 28003735

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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF060515.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P



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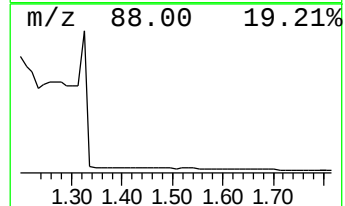
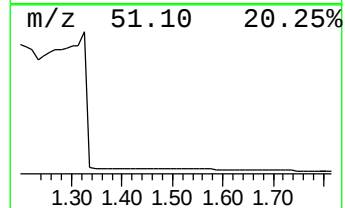
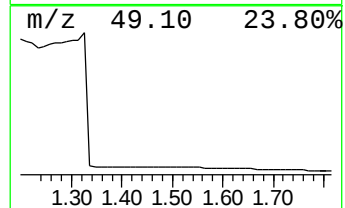
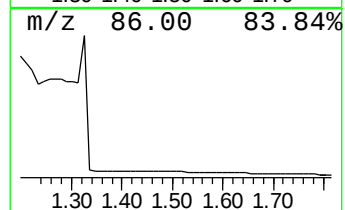
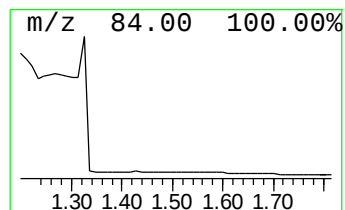
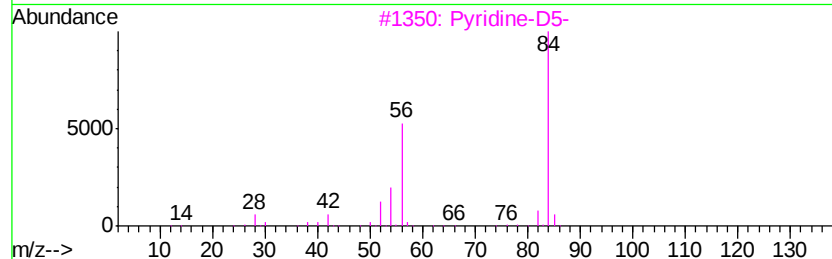
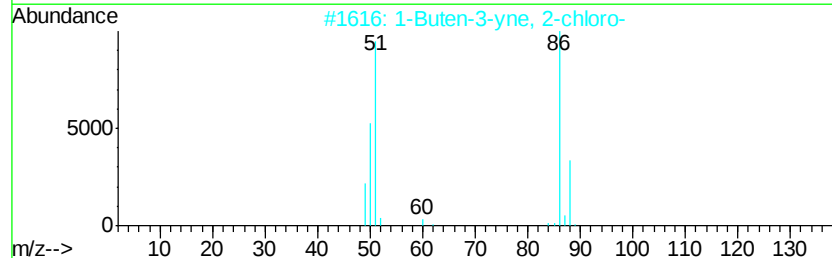
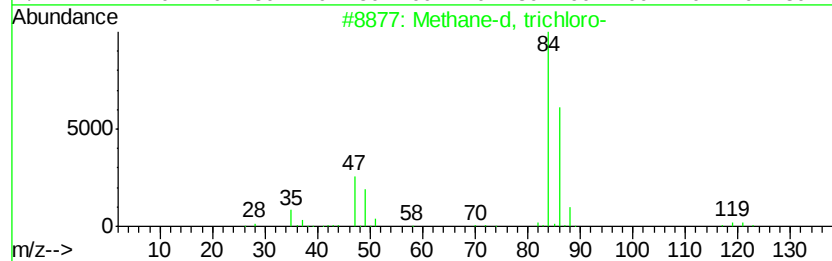
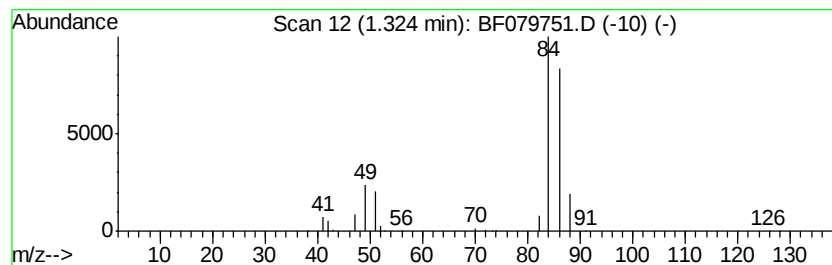
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF060515.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 Methane-d, trichloro- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.32	1518.14 ng	21426000	1,4-Dichlorobenzene-d4	7.48

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Methane-d, trichloro-	119	CDCl3	000865-49-6	9
2		1-Buten-3-yne, 2-chloro-	86	C4H3Cl	017712-36-6	7
3		Pyridine-D5-	84	C5D5N	007291-22-7	5
4		4H-1,2,4-Triazol-4-amine	84	C2H4N4	000584-13-4	5
5		2-Amino-oxazole	84	C3H4N2O	004570-45-0	4



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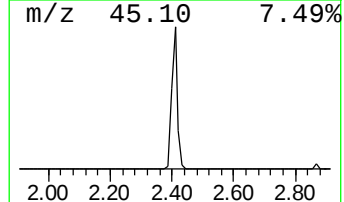
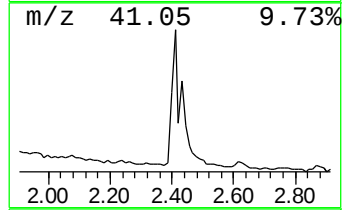
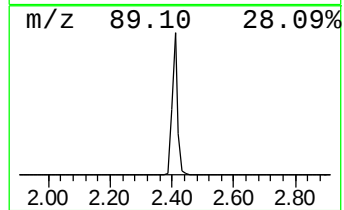
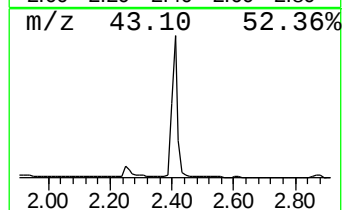
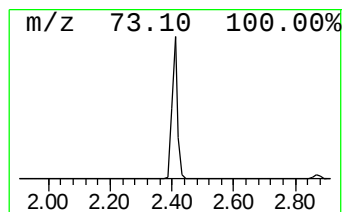
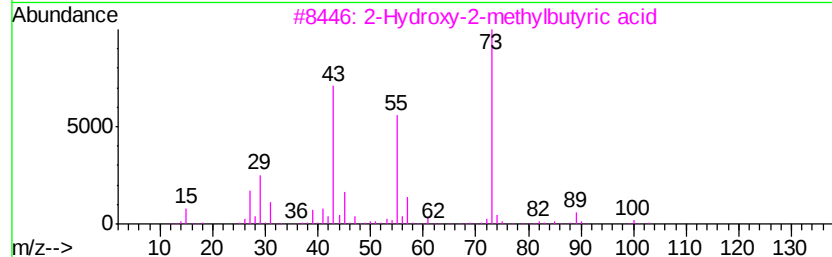
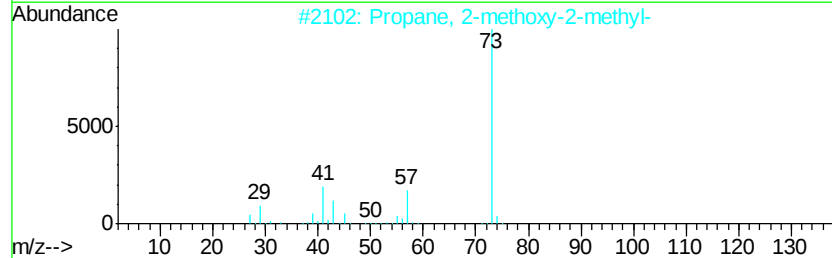
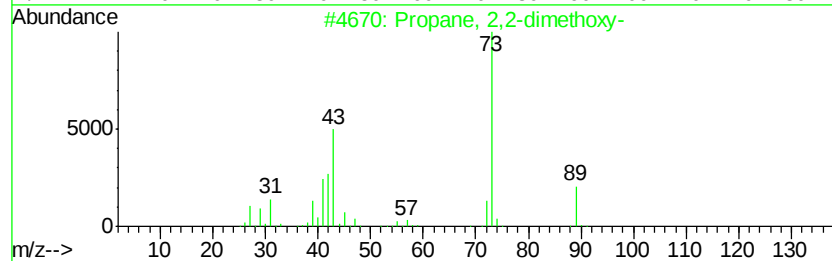
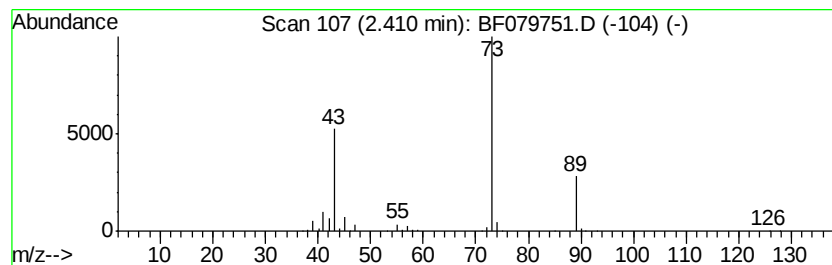
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Peak Number 2 Propane, 2,2-dimethoxy- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.41	46.93 ng	662304	1,4-Dichlorobenzene-d4	7.48

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Propane, 2,2-dimethoxy-	104	C5H12O2	000077-76-9	56
2		Propane, 2-methoxy-2-methyl-	88	C5H12O	001634-04-4	9
3		2-Hydroxy-2-methylbutyric acid	118	C5H10O3	003739-30-8	9
4		1-Hexen-4-ol, 1-chloro-3,5-dimet...	162	C8H15ClO	100244-09-5	9
5		Propanoic acid, 2-methyl-, propy...	130	C7H14O2	000644-49-5	9



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TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Methane-d, trichl...	1.32	1518.1	ng	21426000	1	7.48	282266	20.0
Propane, 2,2-dime...	2.41	46.9	ng	662304	1	7.48	282266	20.0