

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM082718\
Data File : BM016613.D
Acq On : 27 Aug 2018 19:04
Operator : SJ/JU
Sample : J4652-04
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
SONO-W

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:\SVOASRV\HPCHEM1\BNA_M\METHODS\8270-BM082118.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	7.534	712	722	733	rBV	1726097	2903248	9.34%	3.408%
2	8.322	849	856	869	rVB	1758637	2951997	9.50%	3.466%
3	13.263	1690	1696	1704	rVB	5262859	7296920	23.48%	8.567%
4	14.951	1977	1983	1989	rBV3	2406879	5017428	16.14%	5.891%
5	15.286	2031	2040	2048	rBV	22705052	31077482	100.00%	36.485%
6	16.139	2179	2185	2197	rVB	7094181	9999075	32.17%	11.739%
7	19.980	2832	2838	2843	rBV	16938180	18676884	60.10%	21.927%
8	21.533	3097	3102	3112	rVB	2906056	3799428	12.23%	4.461%
9	23.815	3484	3490	3500	rVB	1749155	3455717	11.12%	4.057%

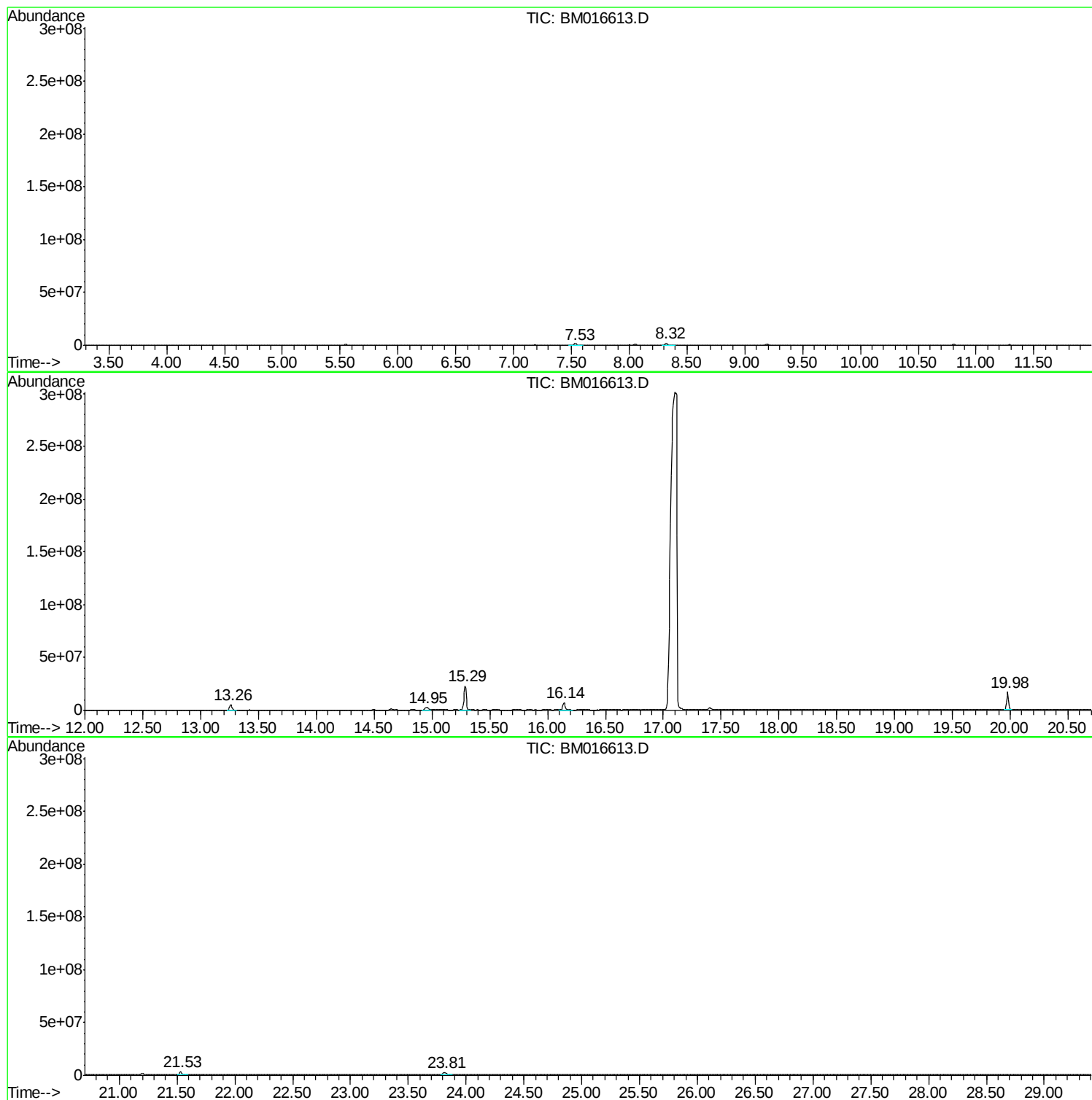
Sum of corrected areas: 85178179

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Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM082118.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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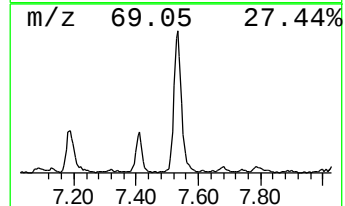
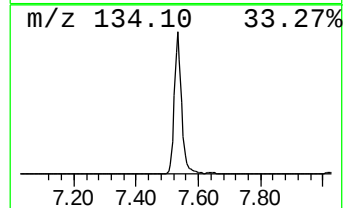
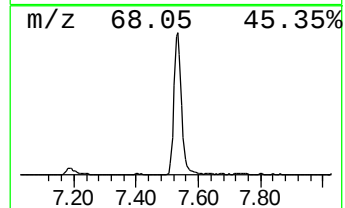
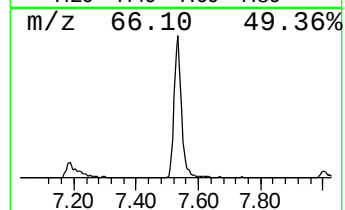
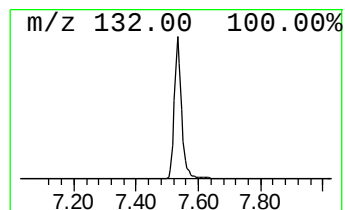
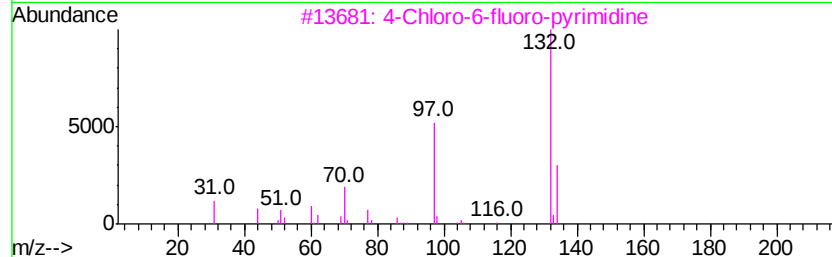
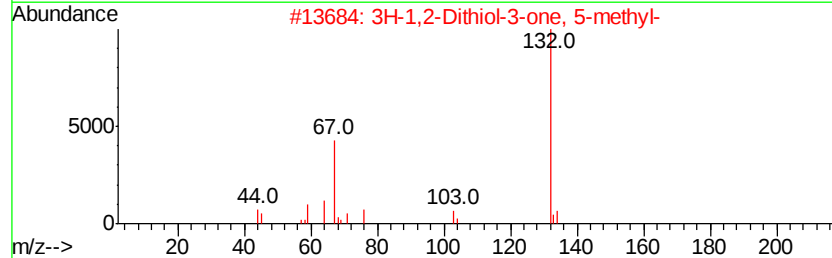
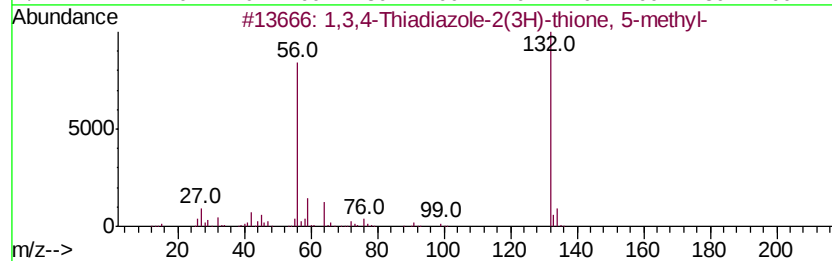
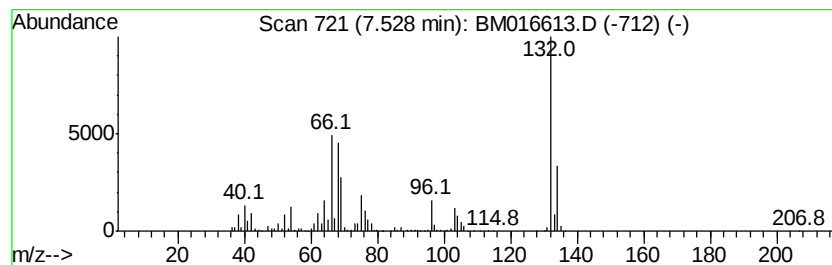
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 unknown7.53 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.53	19.67 ng	2903250	1,4-Dichlorobenzene-d4	8.00

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1,3,4-Thiadiazole-2(3H)-thione, ...	132	C3H4N2S2	029490-19-5	14
2		3H-1,2-Dithiol-3-one, 5-methyl-	132	C4H4OS2	003620-08-4	12
3		4-Chloro-6-fluoro-pyrimidine	132	C4H2ClFN2	051422-01-6	10
4		1,3-Cyclopentanedione, 2-chloro-	132	C5H5ClO2	014203-19-1	10
5		1H-Tetrazaborole, 5-chloro-4,5-d...	132	C2H6BClN4	021960-49-6	9



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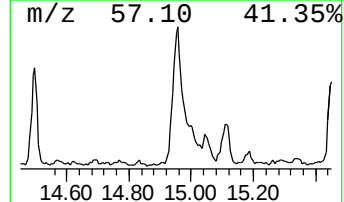
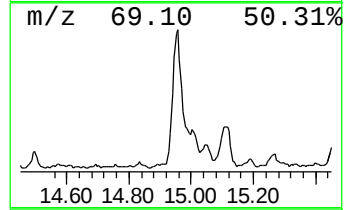
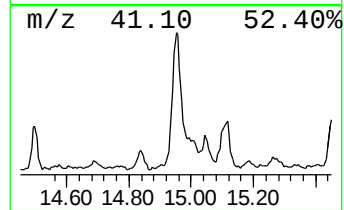
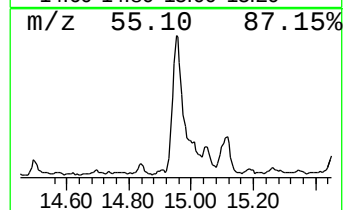
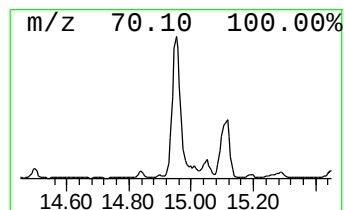
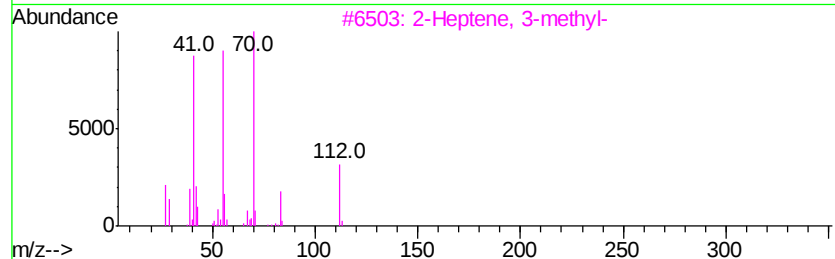
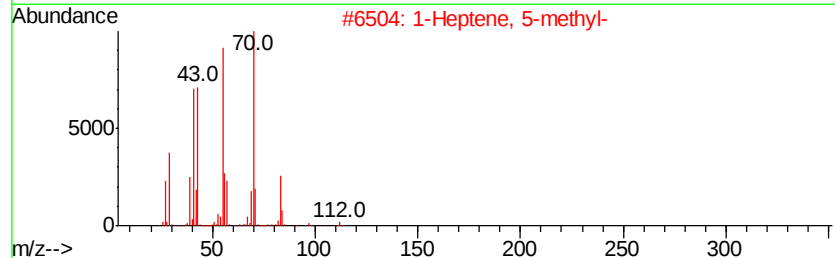
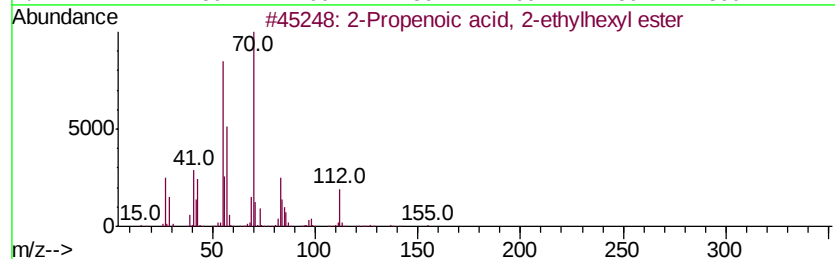
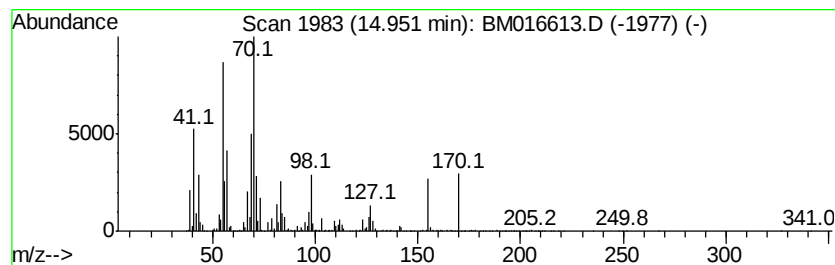
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Peak Number 3 unknown14.95 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.95	20.00 ng	5017430	Acenaphthene-d10	14.65

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Propenoic acid, 2-ethylhexyl e...	184	C11H20O2	000103-11-7	43
2		1-Heptene, 5-methyl-	112	C8H16	013151-04-7	43
3		2-Heptene, 3-methyl-	112	C8H16	003404-75-9	35
4		Cyclopentane, 1,1,3,3-tetramethyl-	126	C9H18	050876-33-0	35
5		Pyrrolidine, 1-(1-oxopropyl)-	127	C7H13NO	004553-05-3	35



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
unknown7.53	7.53	19.7	ng	2903250	1	8.00	2952000	20.0
unknown14.95	14.95	20.0	ng	5017430	3	14.65	5017430	20.0