

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM070918\
Data File : BM015867.D
Acq On : 10 Jul 2018 08:53
Operator : SJ/JU
Sample : J3891-08
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
72-12002-RBR-RBR-VNJ-230

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-BM070518.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	5.316	337	345	356	rBV	225073	352157	4.83%	1.192%
2	5.798	421	427	448	rBV	1261035	1993406	27.34%	6.746%
3	7.445	700	707	725	rBV	1204991	1979205	27.15%	6.698%
4	7.816	764	770	791	rBV	1630214	2630646	36.08%	8.902%
5	8.292	845	851	861	rBV	239367	412003	5.65%	1.394%
6	8.616	900	906	922	rBV	1161619	1923032	26.38%	6.508%
7	9.481	1046	1053	1072	rBV	944209	1590062	21.81%	5.381%
8	11.122	1326	1332	1345	rBV	313185	525940	7.21%	1.780%
9	13.539	1735	1743	1753	rBV	2467180	3591246	49.26%	12.153%
10	14.915	1971	1977	1987	rBV2	459284	702163	9.63%	2.376%
11	16.398	2222	2229	2253	rBV	2449378	3444761	47.25%	11.657%
12	17.645	2435	2441	2453	rBV	649190	940022	12.89%	3.181%
13	20.209	2871	2877	2882	rBV	6144694	7290292	100.00%	24.670%
14	21.756	3134	3140	3146	rBV	860651	1164754	15.98%	3.942%
15	24.150	3541	3547	3557	rBV	505770	1011232	13.87%	3.422%

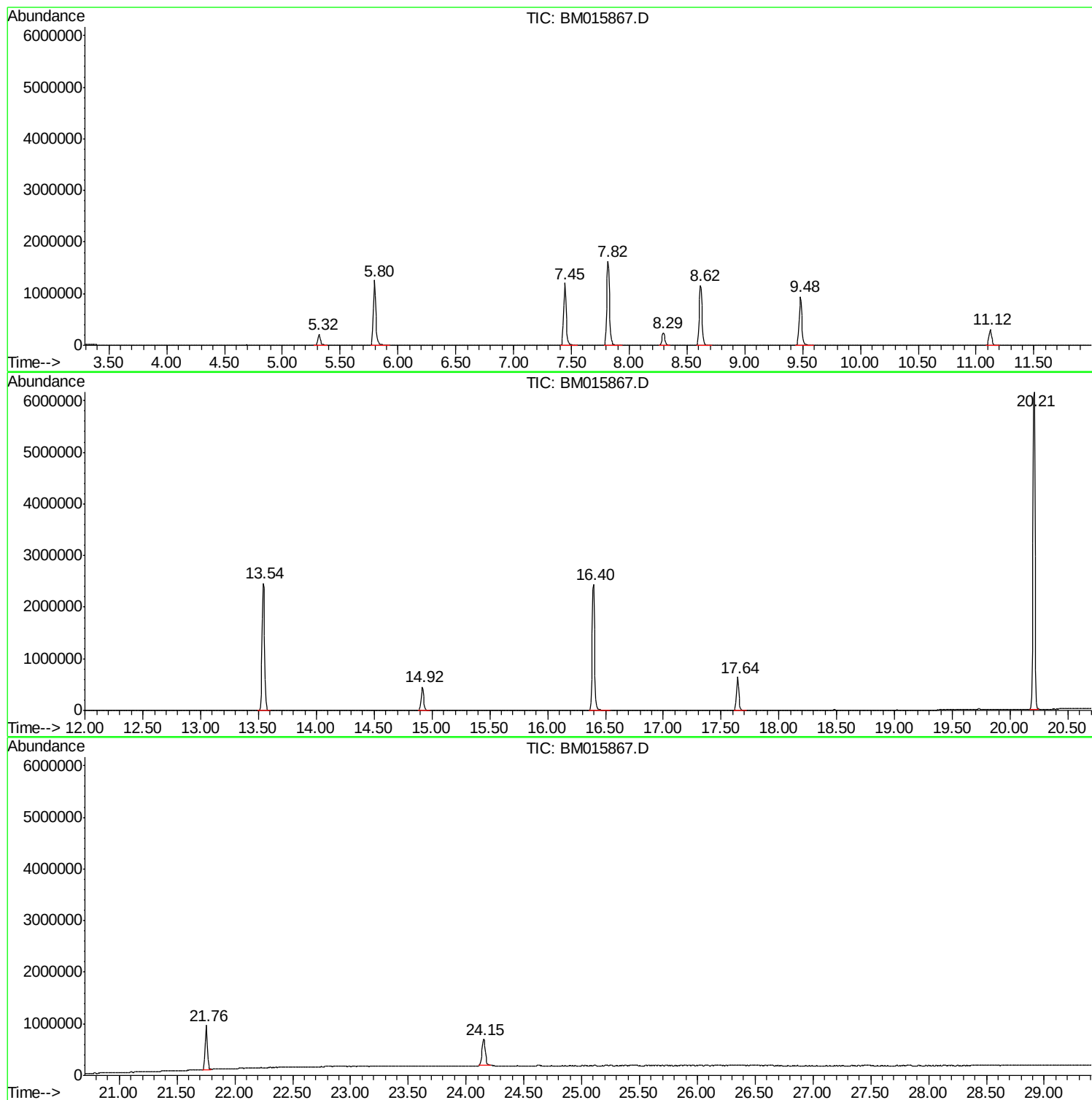
Sum of corrected areas: 29550921

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Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM070518.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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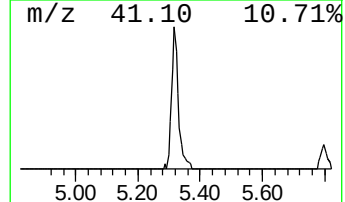
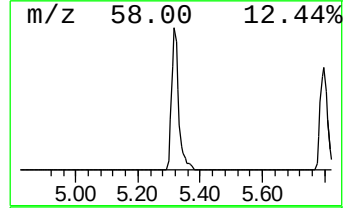
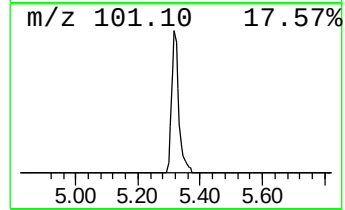
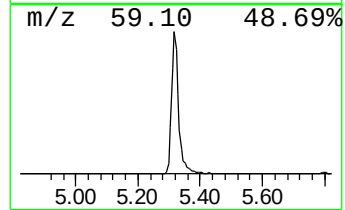
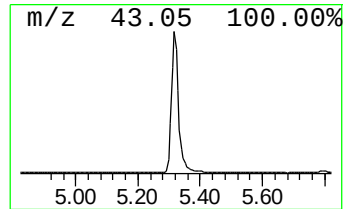
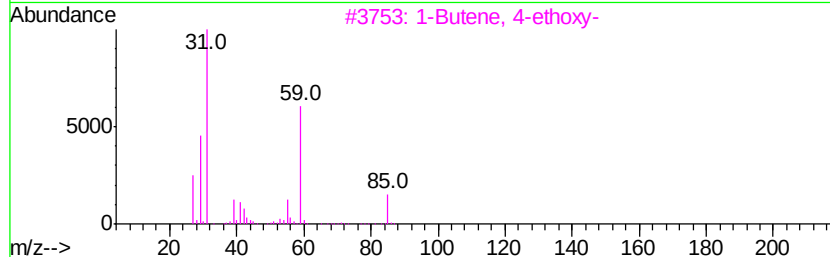
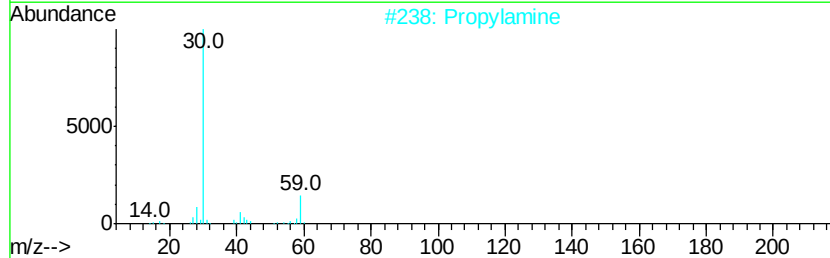
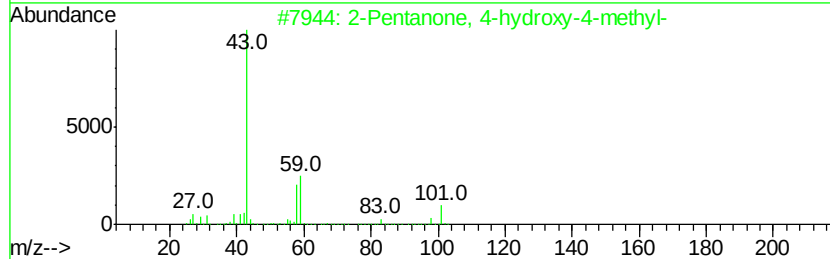
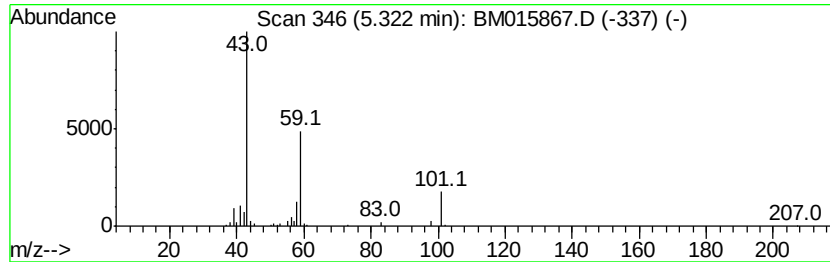
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.32	17.09 ng	352157	1,4-Dichlorobenzene-d4	8.29

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	64
2		Propylamine	59	C3H9N	000107-10-8	40
3		1-Butene, 4-ethoxy-	100	C6H12O	044611-46-3	23
4		2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	16
5		Butane	58	C4H10	000106-97-8	9



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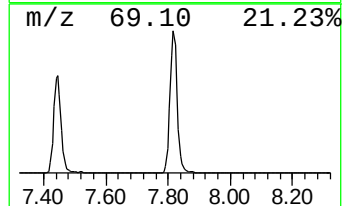
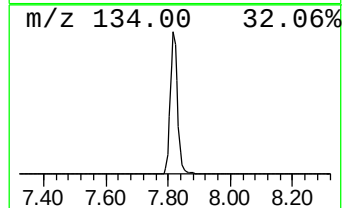
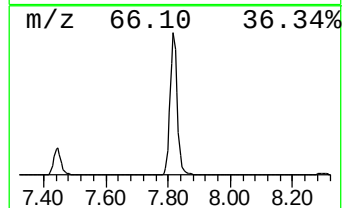
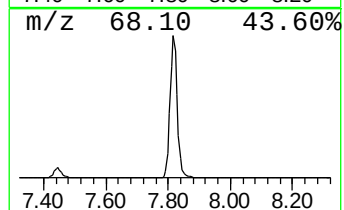
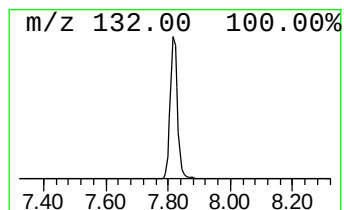
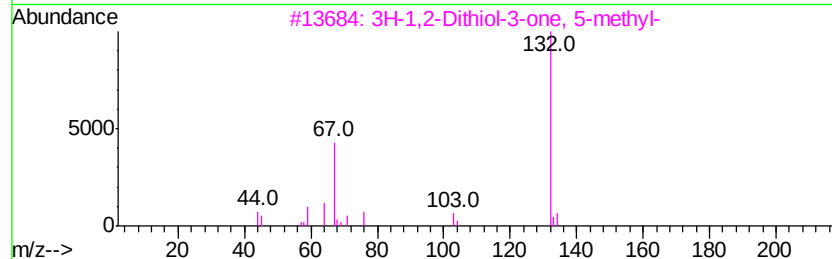
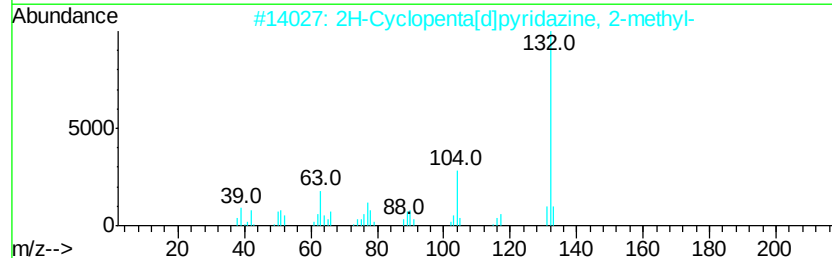
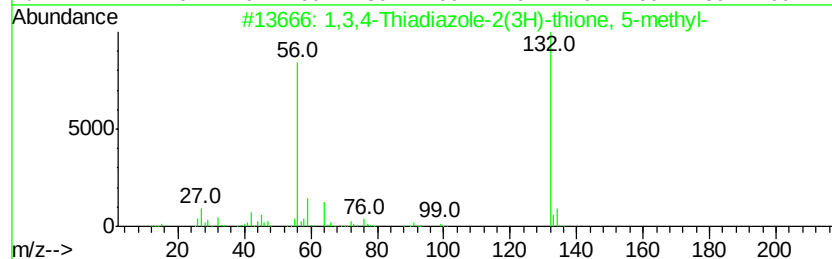
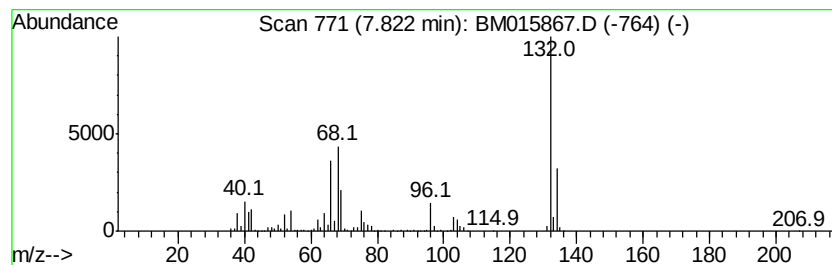
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Peak Number 2 unknown7.82 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.82	127.70 ng	2630650	1,4-Dichlorobenzene-d4	8.29

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1,3,4-Thiadiazole-2(3H)-thione, ...	132	C3H4N2S2	029490-19-5	14
2		2H-Cyclopenta[d]pyridazine, 2-me...	132	C8H8N2	022291-85-6	14
3		3H-1,2-Dithiol-3-one, 5-methyl-	132	C4H4OS2	003620-08-4	14
4		1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	14
5		1,3-Cyclopentanedione, 2-chloro-	132	C5H5ClO2	014203-19-1	12



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TIC Top Hit name	RT	EstConc	Units	Response	---Internal Standard---			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.32	17.1	ng	352157	1	8.29	412003	20.0
unknown7.82	7.82	127.7	ng	2630650	1	8.29	412003	20.0