

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM070518\
 Data File : BM015789.D
 Acq On : 06 Jul 2018 07:05
 Operator : SJ/JU
 Sample : PB110790BL
 Misc :
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 PB110790BL

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	5.328	342	347	360	rBV	351665	553797	5.77%	1.027%
2	5.816	423	430	450	rBV	3457704	5150469	53.66%	9.554%
3	7.469	703	711	726	rBV	3292894	5497524	57.27%	10.198%
4	7.840	766	774	786	rBV	3583925	5820404	60.64%	10.797%
5	8.310	848	854	863	rBV	491593	800174	8.34%	1.484%
6	8.634	901	909	920	rBV	2587553	4195922	43.71%	7.783%
7	9.498	1048	1056	1071	rBV	1985006	3321130	34.60%	6.161%
8	11.139	1328	1335	1347	rBV	569506	984213	10.25%	1.826%
9	13.557	1739	1746	1760	rBV	5367418	7356515	76.64%	13.646%
10	14.927	1973	1979	1989	rBV2	802506	1222434	12.74%	2.268%
11	16.410	2224	2231	2246	rBV	4374794	6020433	62.72%	11.168%
12	17.657	2436	2443	2453	rBV2	945463	1383087	14.41%	2.566%
13	20.215	2872	2878	2884	rBV	8117058	9598714	100.00%	17.805%
14	21.762	3136	3141	3148	rBV	884655	1159433	12.08%	2.151%
15	24.168	3544	3550	3558	rVB	431600	845355	8.81%	1.568%

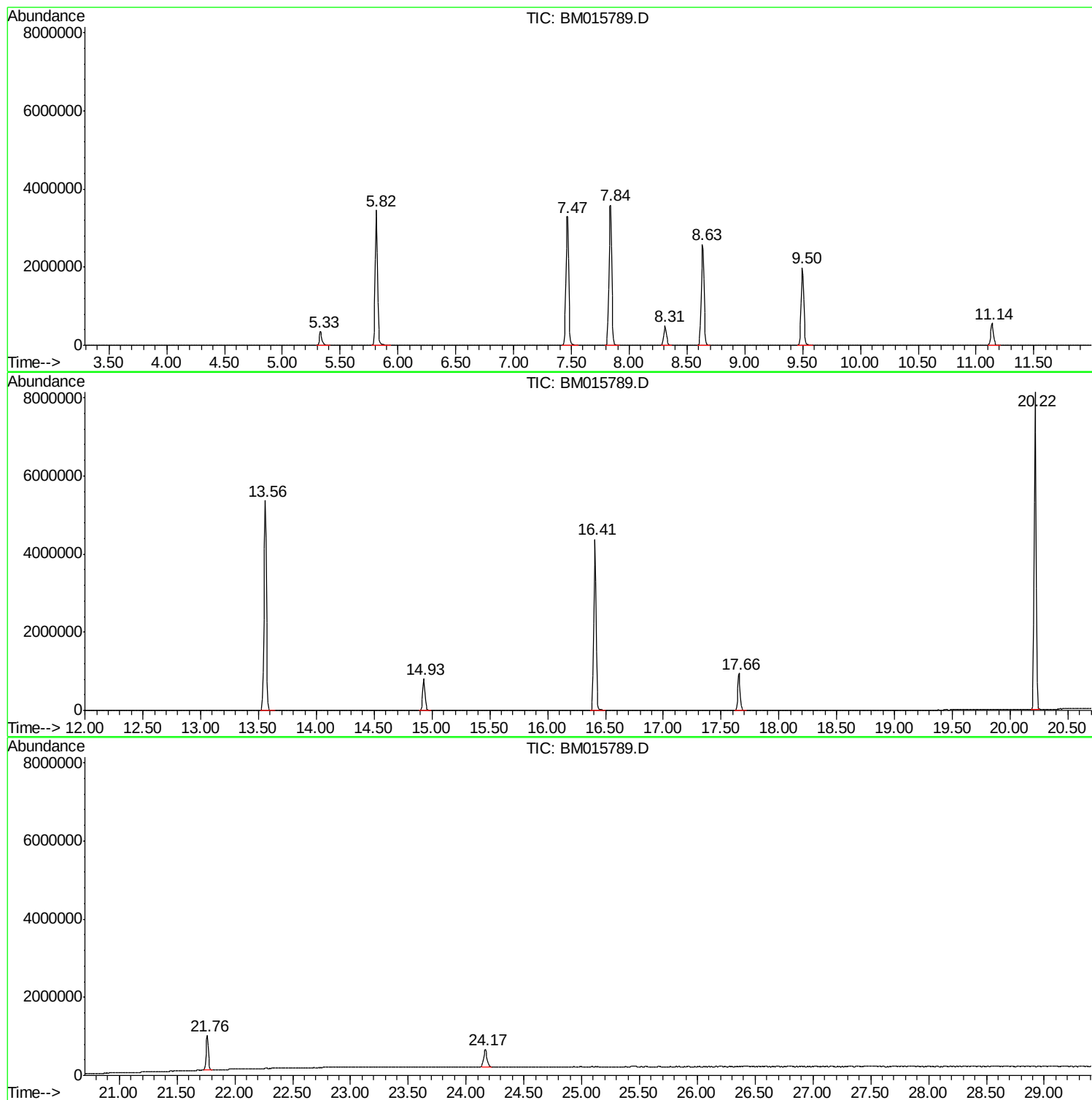
Sum of corrected areas: 53909604

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



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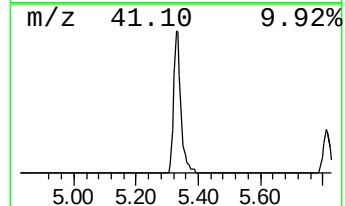
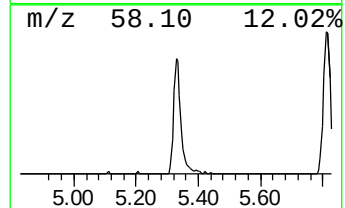
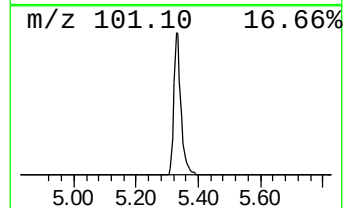
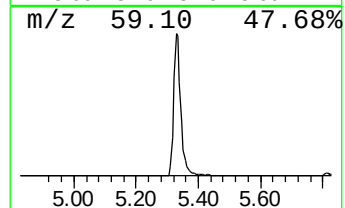
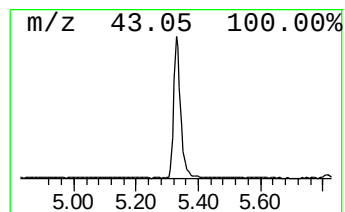
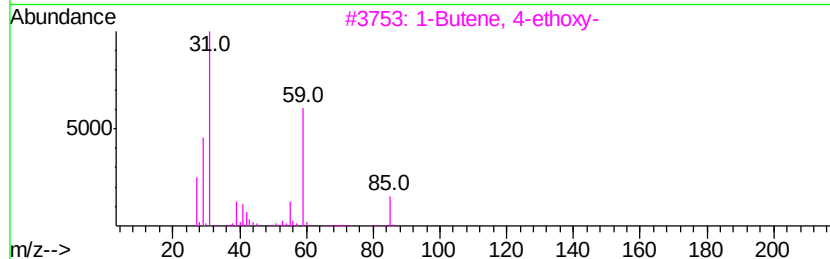
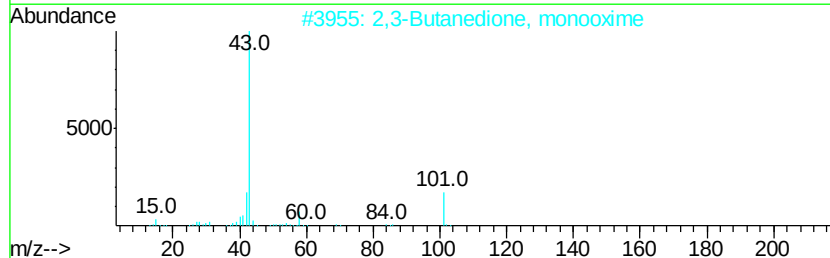
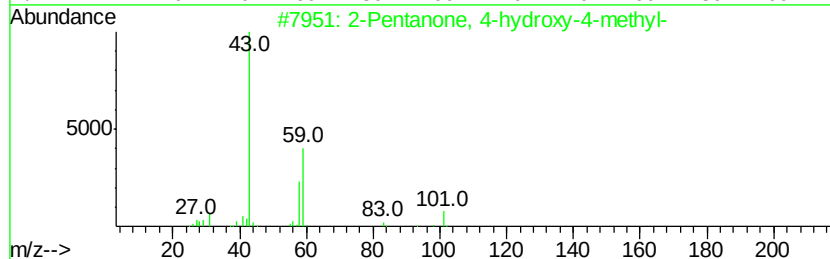
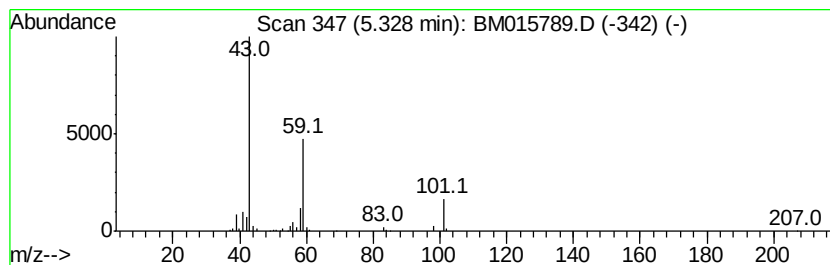
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.33	13.84 ng	553797	1,4-Dichlorobenzene-d4	8.31

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50
2		2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	27
3		1-Butene, 4-ethoxy-	100	C6H12O	044611-46-3	23
4		Butane	58	C4H10	000106-97-8	9
5		Acetic acid, 2-propenyl ester	100	C5H8O2	000591-87-7	9



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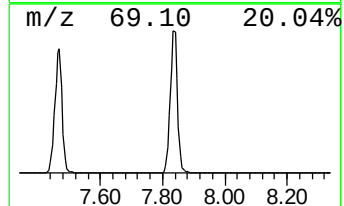
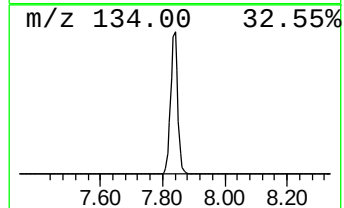
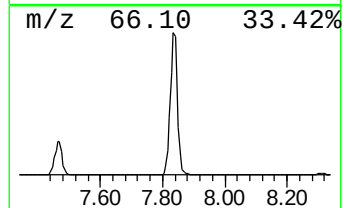
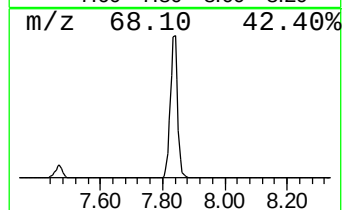
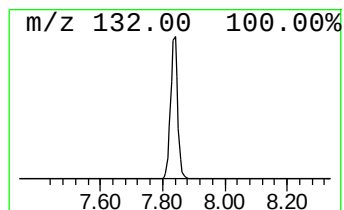
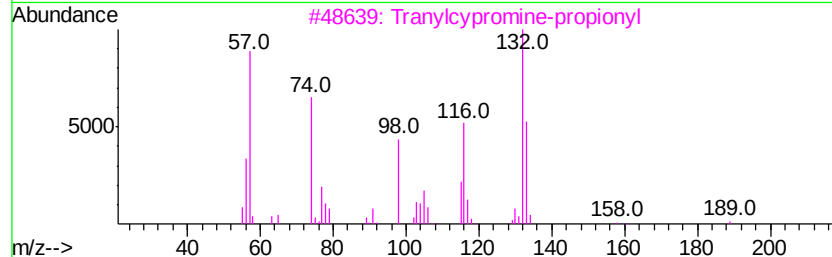
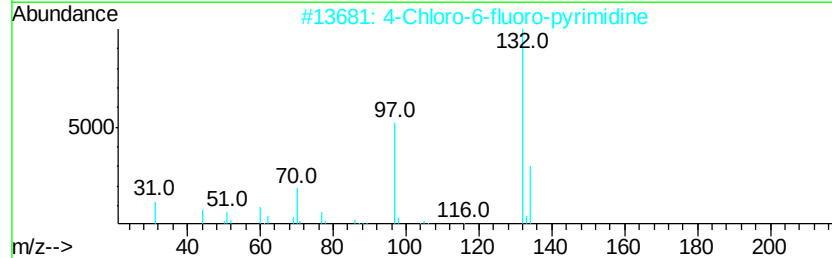
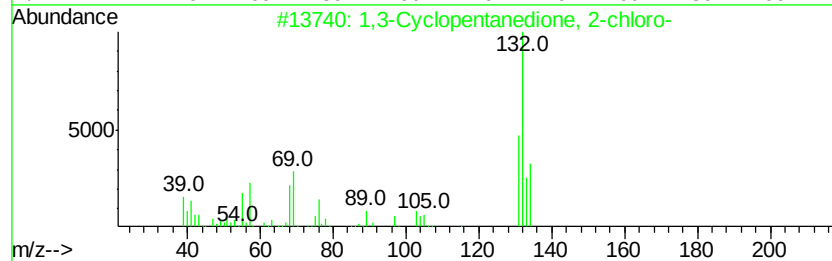
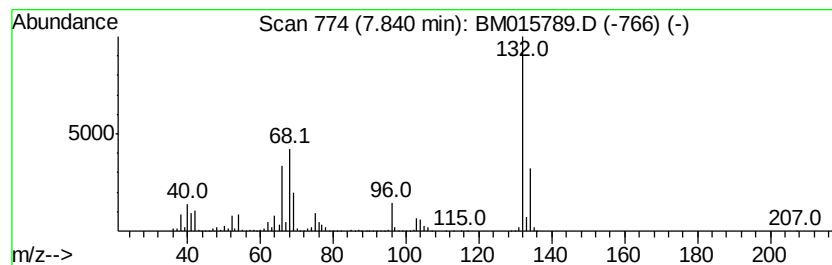
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Peak Number 2 unknown7.84 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.84	145.48 ng	5820400	1,4-Dichlorobenzene-d4	8.31

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1,3-Cyclopentanedione, 2-chloro-	132	C5H5ClO2	014203-19-1	35
2		4-Chloro-6-fluoro-pyrimidine	132	C4H2ClFN2	051422-01-6	27
3		Tranvlcypromine-propionyl	189	C12H15NO	1000123-86-3	22
4		1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	22
5		Imidazole, 4,5-dicyano-1-methyl-	132	C6H4N4	019485-35-9	22



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TIC Top Hit name	RT	EstConc	Units	Response	---Internal Standard---			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.33	13.8	ng	553797	1	8.31	800174	20.0
unknown7.84	7.84	145.5	ng	5820400	1	8.31	800174	20.0