

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM111418\
 Data File : BM017661.D
 Acq On : 14 Nov 2018 23:28
 Operator : MJ/SJ
 Sample : PB114805BL
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 PB114805BL

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-BM111218.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	4.793	301	307	316	rBV	170730	241415	2.09%	0.341%
2	5.240	376	383	399	rBV	4570987	6369183	55.25%	8.997%
3	6.804	641	649	664	rBV	4286177	6756180	58.61%	9.543%
4	7.151	701	708	726	rBV	4508730	6943930	60.24%	9.809%
5	7.610	779	786	794	rBV	928290	1469659	12.75%	2.076%
6	7.928	832	840	851	rBV	3234320	5188960	45.02%	7.330%
7	8.769	972	983	1000	rBV	2363505	4136126	35.88%	5.843%
8	10.381	1250	1257	1279	rVB	1003624	1966438	17.06%	2.778%
9	12.869	1673	1680	1697	rBV	6229511	9420625	81.73%	13.307%
10	14.257	1909	1916	1930	rBV2	1543142	2528216	21.93%	3.571%
11	15.757	2164	2171	2194	rBV	4422909	6783147	58.85%	9.582%
12	17.010	2377	2384	2408	rBV2	1698202	2887166	25.05%	4.078%
13	19.651	2827	2833	2846	rBV	9065773	11527007	100.00%	16.283%
14	21.209	3093	3098	3110	rBV	1738673	2632709	22.84%	3.719%
15	23.403	3465	3471	3484	rBV	938813	1942935	16.86%	2.745%

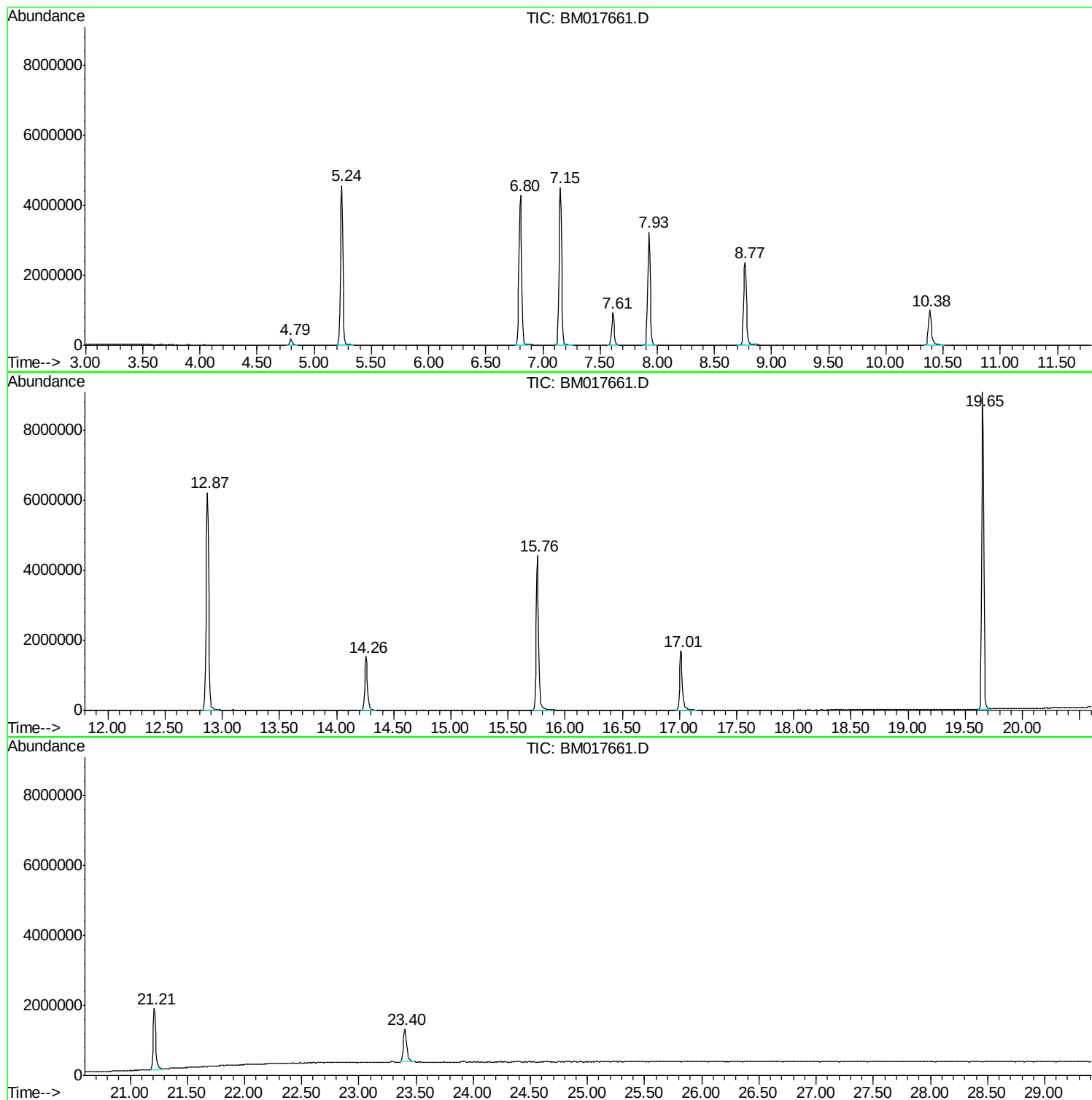
Sum of corrected areas: 70793696

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Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM111218.M
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P



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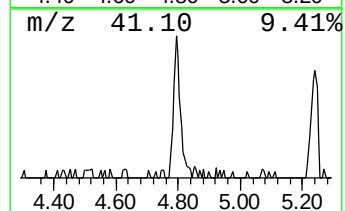
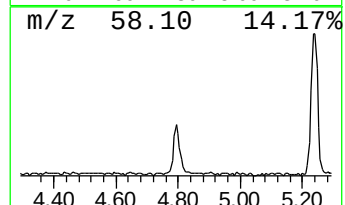
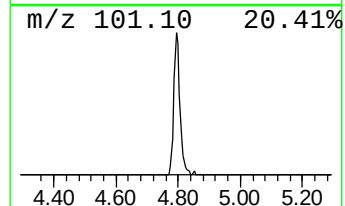
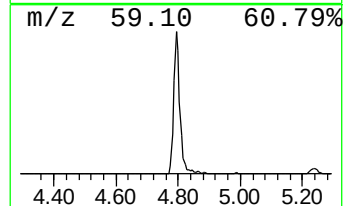
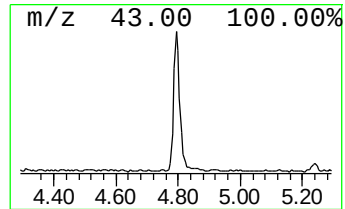
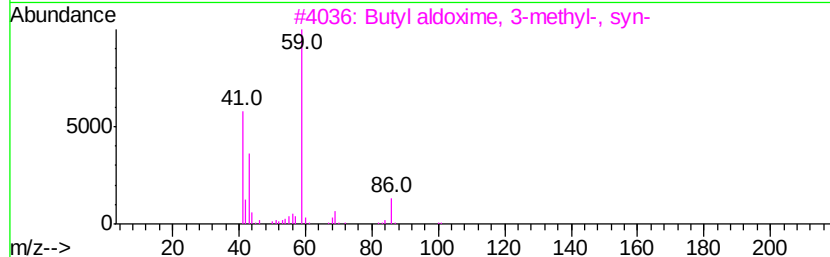
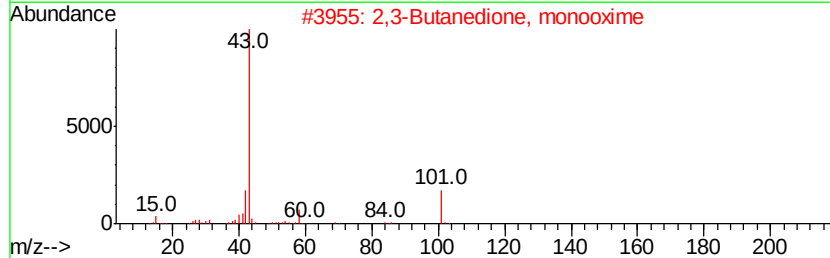
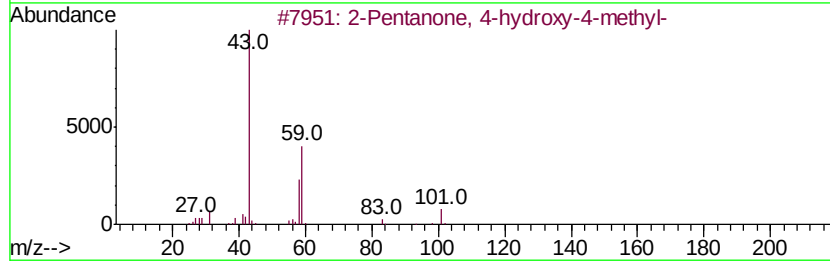
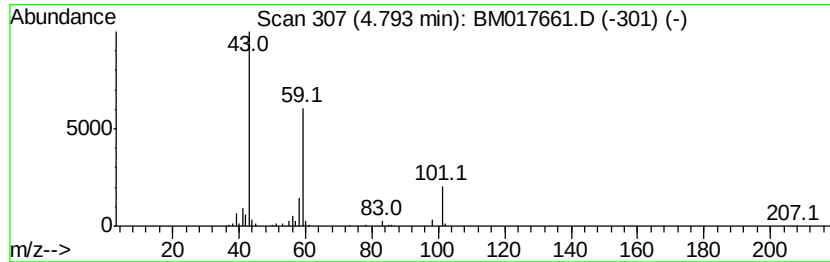
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.79	3.29 ng	241415	1,4-Dichlorobenzene-d4	7.61

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56
2		2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	35
3		Butyl aldoxime, 3-methyl-, syn-	101	C5H11NO	005780-40-5	25
4		1-Butene, 4-ethoxy-	100	C6H12O	044611-46-3	17
5		Butane	58	C4H10	000106-97-8	9



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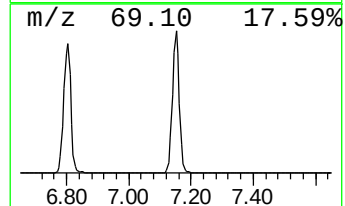
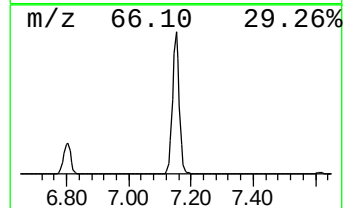
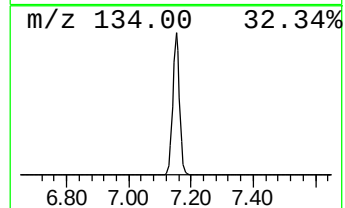
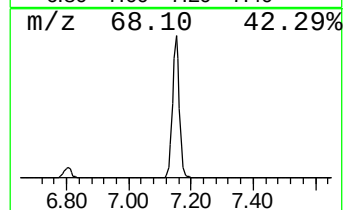
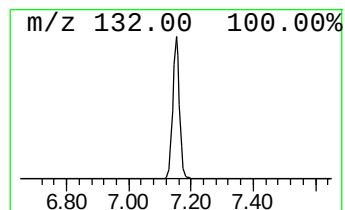
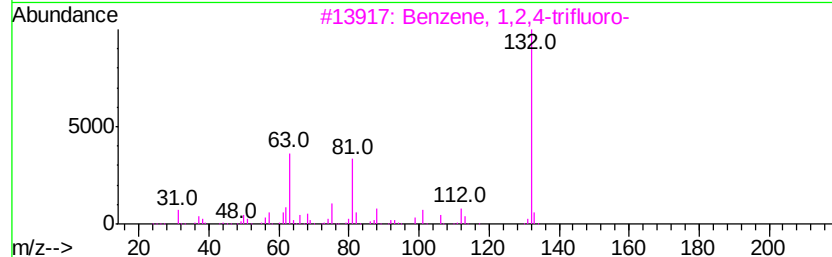
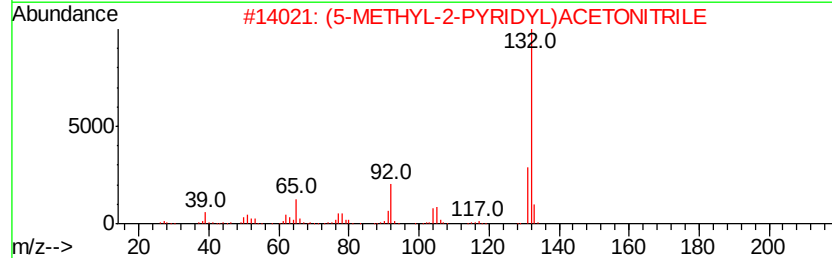
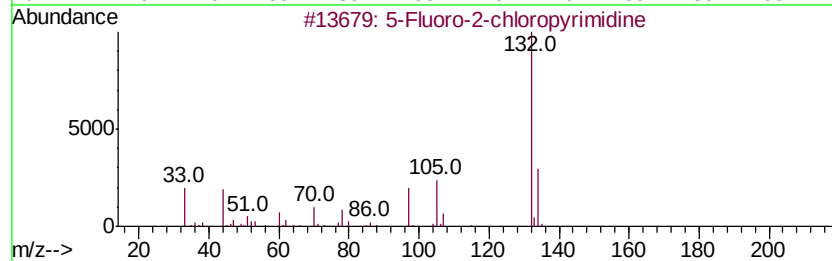
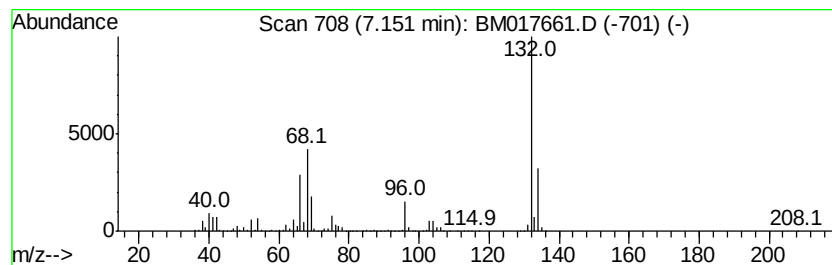
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Peak Number 2 unknown7.15 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.15	94.50 ng	6943930	1,4-Dichlorobenzene-d4	7.61

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	12
2		(5-METHYL-2-PYRIDYL)ACETONITRILE	132	C8H8N2	1000241-93-9	10
3		Benzene, 1,2,4-trifluoro-	132	C6H3F3	000367-23-7	9
4		Benzene, 1,3,5-trifluoro-	132	C6H3F3	000372-38-3	9
5		Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9



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TIC Top Hit name	RT	EstConc	Units	Response	---Internal Standard---			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	4.79	3.3	ng	241415	1	7.61	1469660	20.0
unknown7.15	7.15	94.5	ng	6943930	1	7.61	1469660	20.0