

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM071218\
 Data File : BM015924.D
 Acq On : 12 Jul 2018 20:44
 Operator : SJ/JU
 Sample : PB111042BL
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 PB111042BL

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	5.310	339	344	358	rBV	76146	130490	2.80%	0.539%
2	5.793	420	426	446	rBV	1150304	1802797	38.66%	7.444%
3	7.440	699	706	721	rBV	1133840	1881605	40.35%	7.770%
4	7.810	763	769	785	rBV	1249709	2031841	43.57%	8.390%
5	8.287	844	850	862	rVB	316105	517144	11.09%	2.135%
6	8.610	898	905	921	rBV	1435198	2303665	49.40%	9.512%
7	9.475	1045	1052	1067	rBV	1049829	1780084	38.17%	7.350%
8	11.116	1324	1331	1352	rBV	350740	628178	13.47%	2.594%
9	13.533	1735	1742	1757	rBV	2643244	3768659	80.81%	15.562%
10	14.910	1970	1976	1987	rBV2	473009	723579	15.52%	2.988%
11	16.386	2221	2227	2243	rBV	1242756	1719417	36.87%	7.100%
12	17.633	2434	2439	2449	rBV2	529694	775985	16.64%	3.204%
13	20.198	2869	2875	2882	rBV	4204600	4663453	100.00%	19.257%
14	21.745	3133	3138	3146	rBV	583465	772692	16.57%	3.191%
15	24.139	3539	3545	3555	rVB	341732	717799	15.39%	2.964%

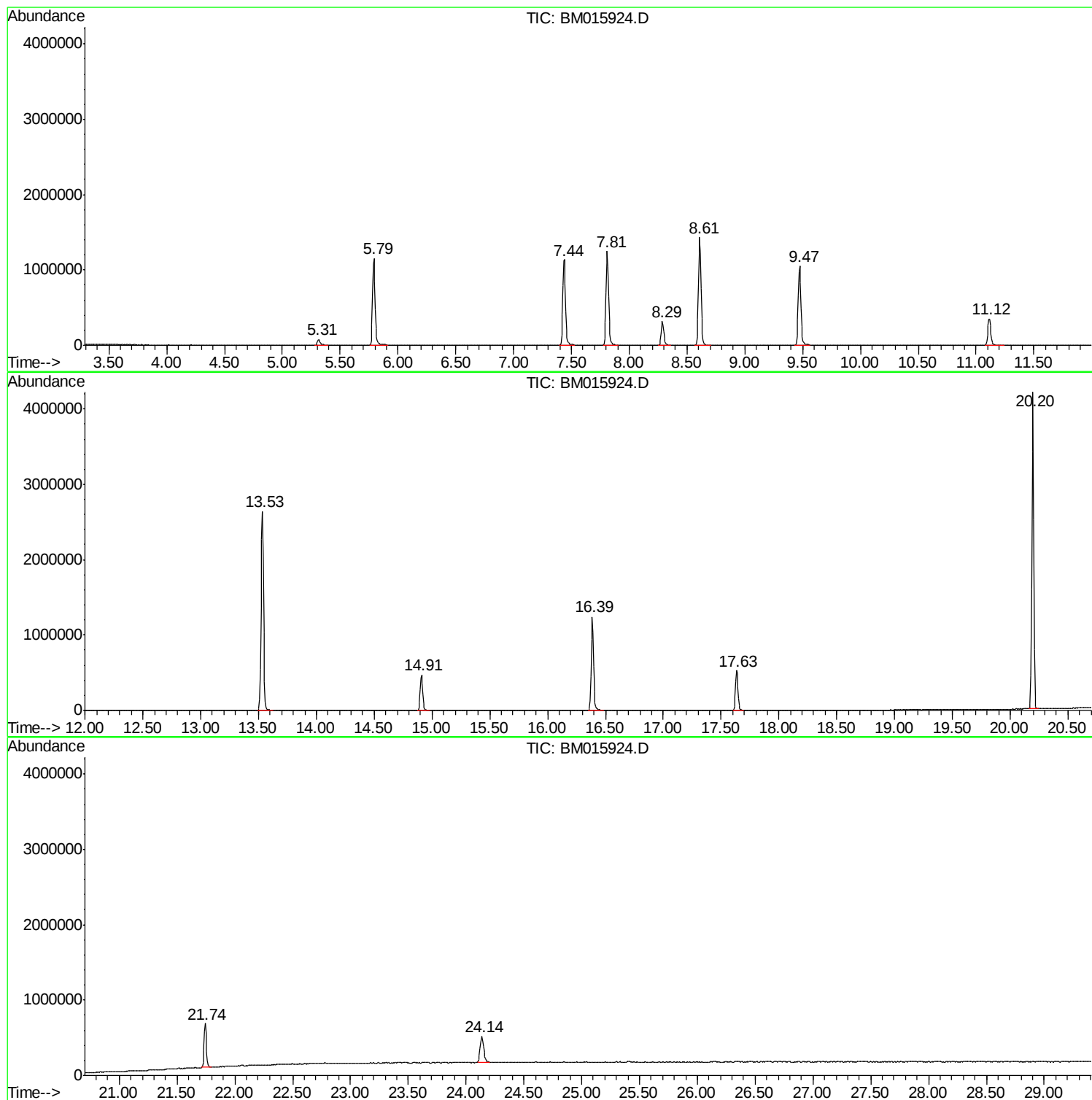
Sum of corrected areas: 24217388

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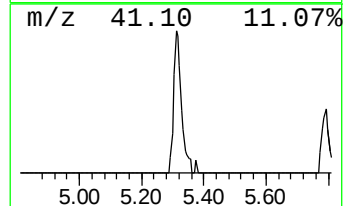
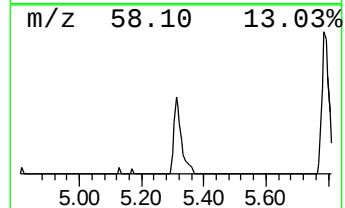
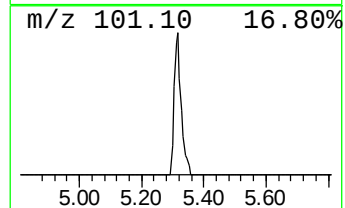
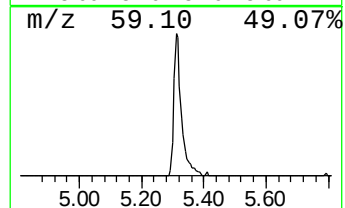
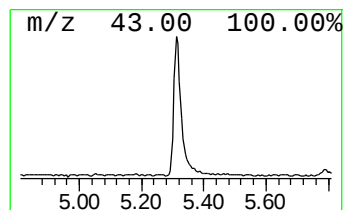
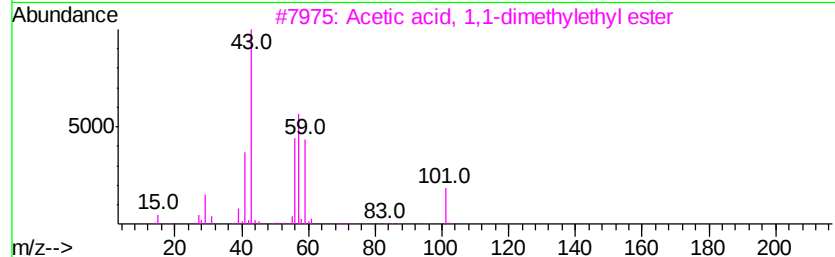
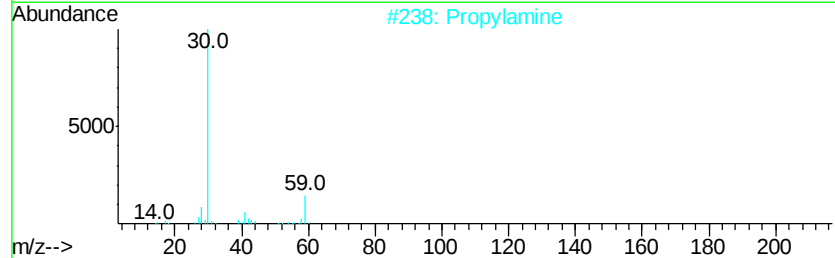
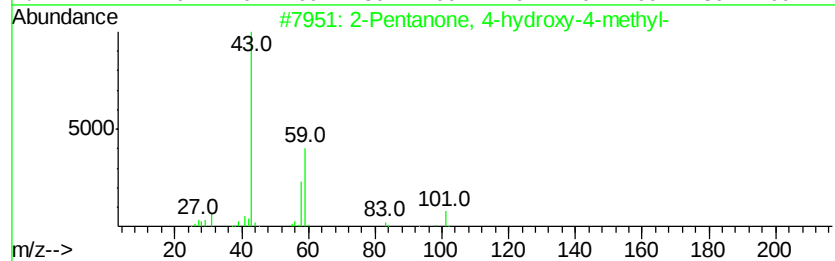
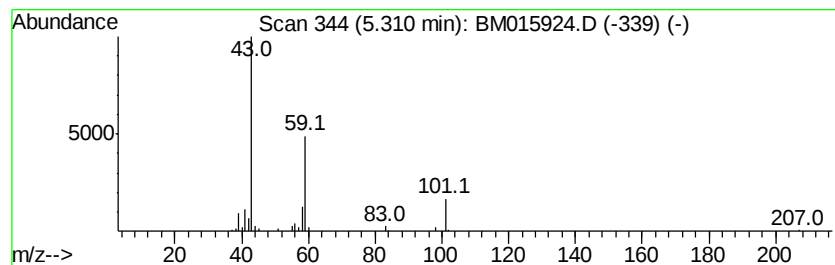
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.31	5.05 ng	130490	1,4-Dichlorobenzene-d4	8.29

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50
2			Propylamine	59	C3H9N	000107-10-8	40
3			Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	38
4			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	27
5			1-Butene, 4-ethoxy-	100	C6H12O	044611-46-3	23



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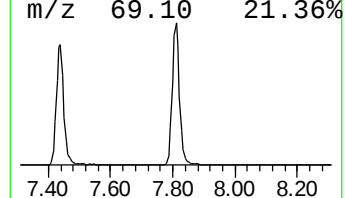
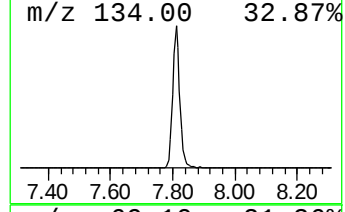
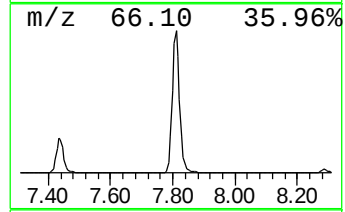
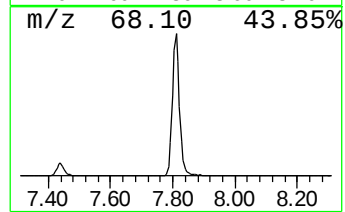
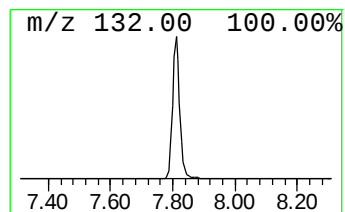
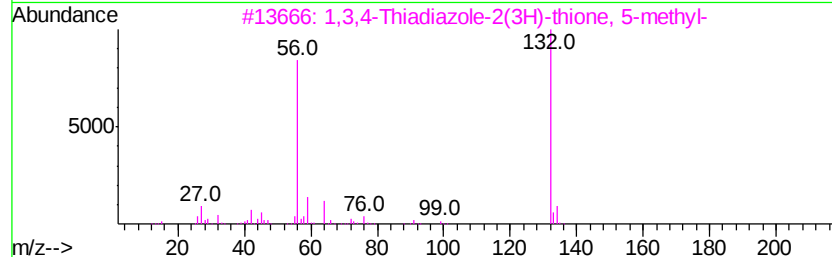
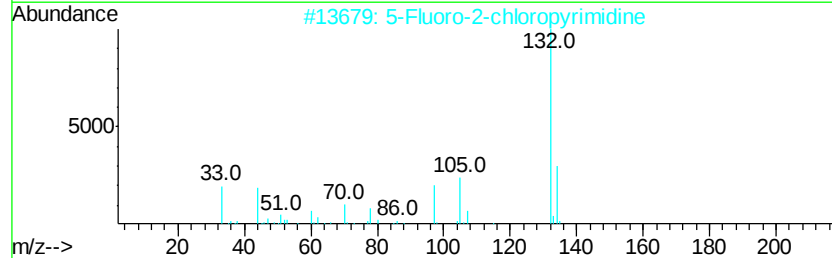
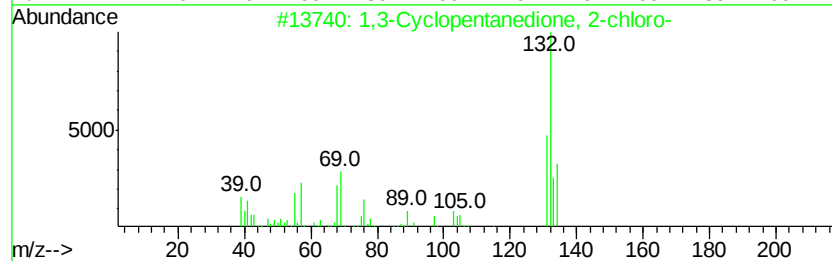
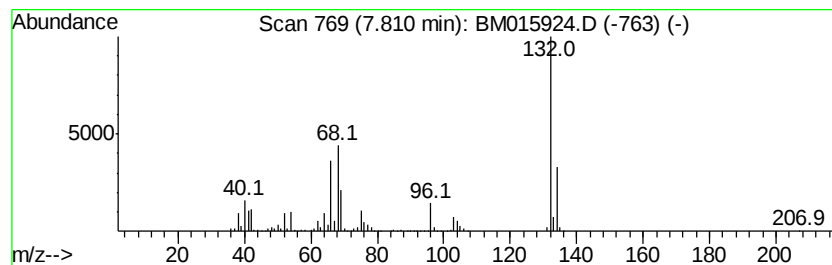
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Peak Number 2 unknown7.81 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.81	78.58 ng	2031840	1,4-Dichlorobenzene-d4	8.29

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1,3-Cyclopentanedione, 2-chloro-	132	C5H5ClO2	014203-19-1	35
2		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	27
3		1,3,4-Thiadiazole-2(3H)-thione, ...	132	C3H4N2S2	029490-19-5	27
4		1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	27
5		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	22



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.31	5.0	ng	130490	1	8.29	517144	20.0
unknown7.81	7.81	78.6	ng	2031840	1	8.29	517144	20.0