

Data Path : Z:\HPCHEM1\BNA N\DATA\BN040918\
 Data File : BN000728.D
 Acq On : 10 Apr 2018 02:01
 Operator : JU/SJ
 Sample : J2292-05
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 040618-JETW-E-U-M

Integration Parameters: rteint.p
 Integrator: RTE
 Smoothing : ON
 Sampling : 1
 Start Thrs: 0.2
 Stop Thrs : 0
 Filtering: 5
 Min Area: 3 % of largest Peak
 Max Peaks: 100
 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
 Peak separation: 5

Method : Z:\HPCHEM1\BNA_N\METHODS\8270-BN040918.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	4.811	305	310	320	rBV	81555	110068	1.63%	0.307%
2	5.258	380	386	397	rBV	851291	1158548	17.12%	3.231%
3	6.811	644	650	662	rVB	478720	724032	10.70%	2.019%
4	7.169	704	711	720	rBV	1634279	2496832	36.90%	6.962%
5	7.634	783	790	798	rVB	586526	900025	13.30%	2.510%
6	7.952	836	844	852	rVB	2437161	3843035	56.79%	10.716%
7	8.781	976	985	992	rBV	1841229	2985813	44.13%	8.326%
8	10.399	1253	1260	1267	rBV	788812	1259389	18.61%	3.512%
9	12.887	1675	1683	1689	rBV	4492965	6596767	97.49%	18.395%
10	13.728	1821	1826	1831	rBV	98155	125719	1.86%	0.351%
11	14.263	1911	1917	1929	rBV2	1019653	1519669	22.46%	4.237%
12	15.151	2063	2068	2073	rVB	74781	90017	1.33%	0.251%
13	15.751	2165	2170	2186	rVB	1849617	2528500	37.37%	7.050%
14	16.016	2210	2215	2227	rBV	61273	99185	1.47%	0.277%
15	17.010	2378	2384	2393	rVB	1132518	1555939	22.99%	4.339%
16	17.928	2536	2540	2548	rBV	51206	72417	1.07%	0.202%
17	19.657	2828	2834	2840	rBV	5205248	6766665	100.00%	18.868%
18	20.957	3051	3055	3066	rBV	62323	111655	1.65%	0.311%
19	21.210	3093	3098	3105	rBV	1180434	1495670	22.10%	4.171%
20	23.404	3464	3471	3481	rVB	818662	1422764	21.03%	3.967%

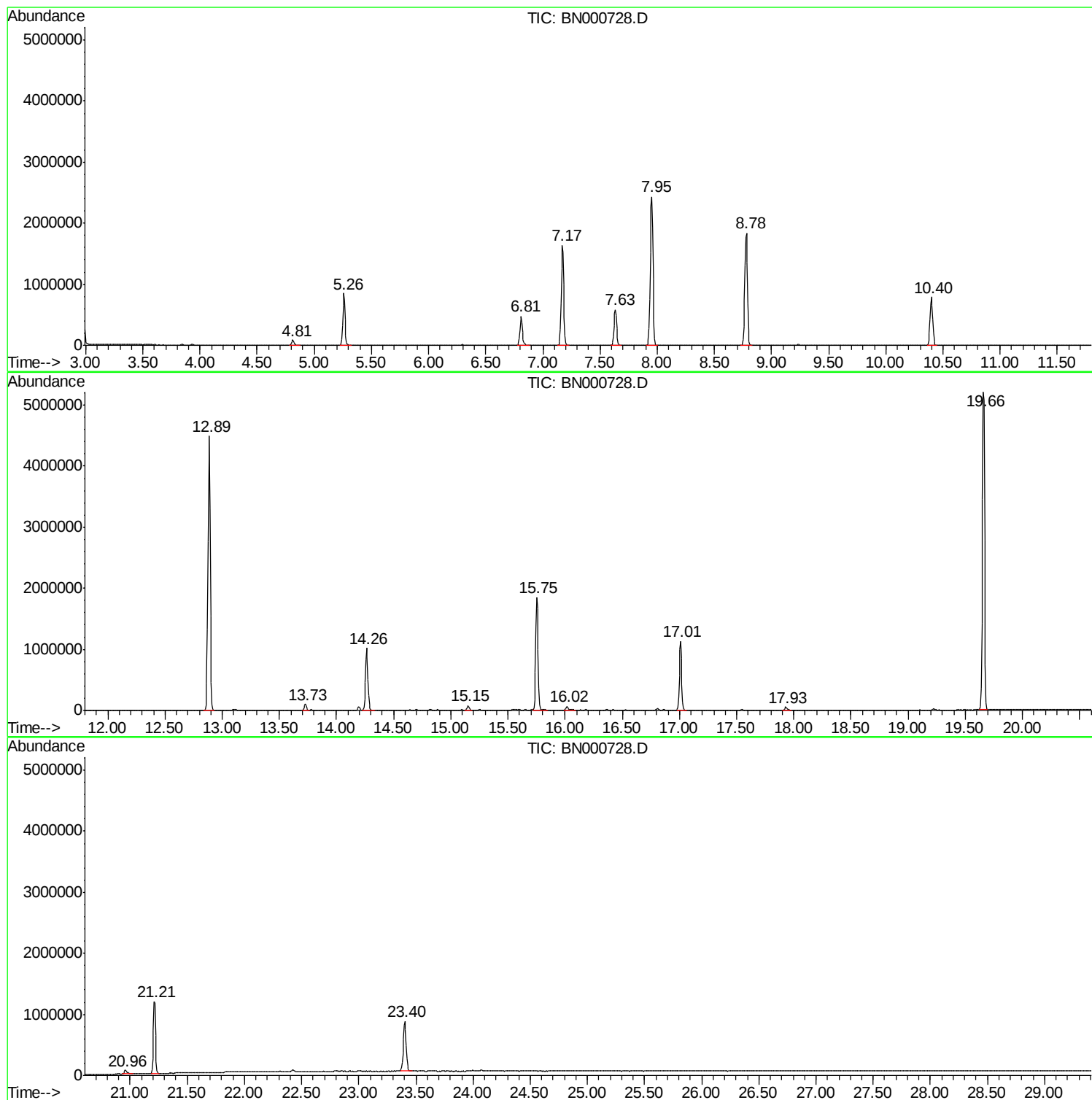
Sum of corrected areas: 35862709

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Quant Method : Z:\HPCHEM1\BNA N\METHODS\8270-BN040918.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P



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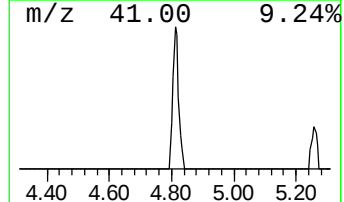
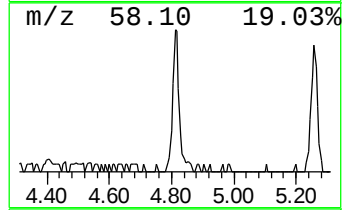
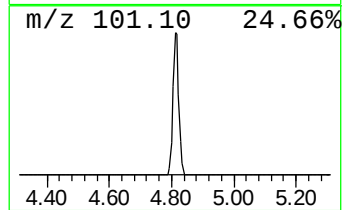
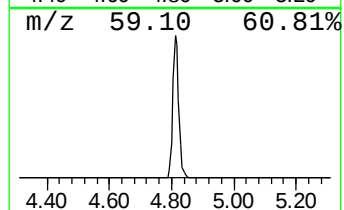
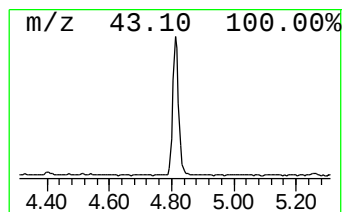
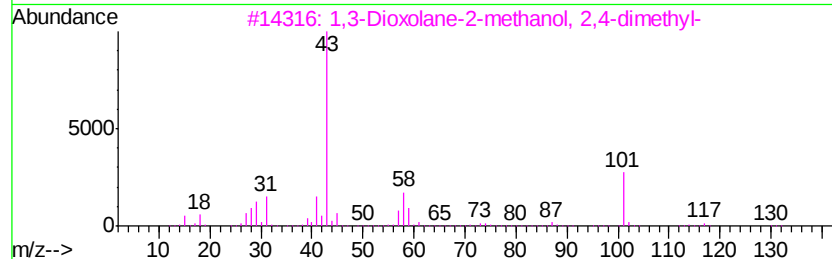
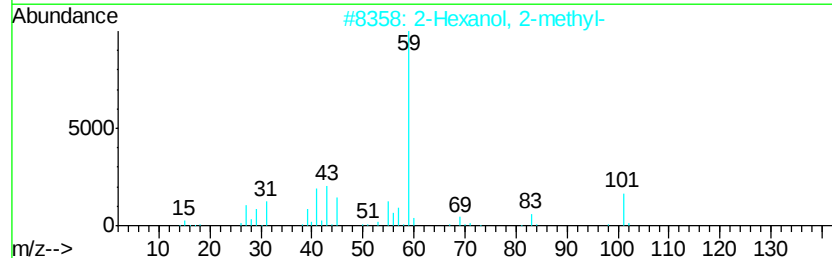
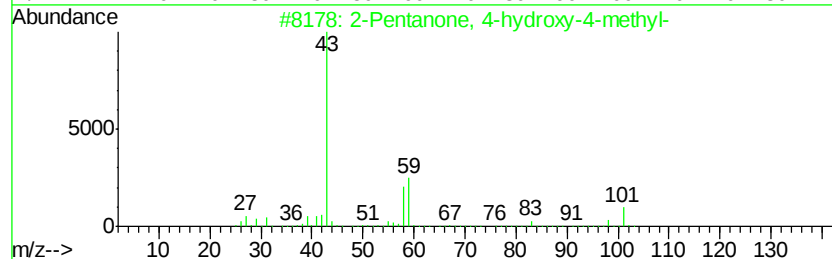
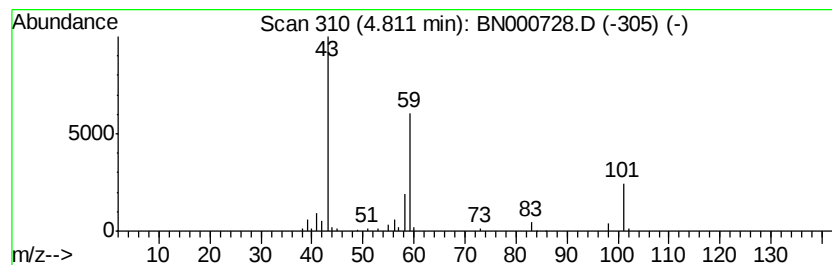
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.81	2.45 ng	110068	1,4-Dichlorobenzene-d4	7.63

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	53
2			2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	25
3			1,3-Dioxolane-2-methanol, 2,4-di...	132	C6H12O3	053951-43-2	25
4			Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	12
5			Acetone	58	C3H6O	000067-64-1	10



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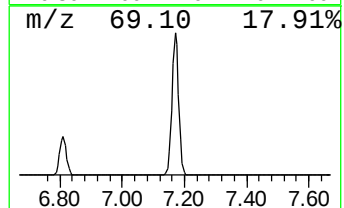
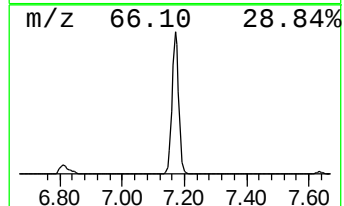
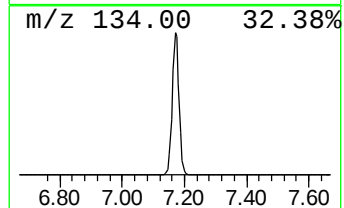
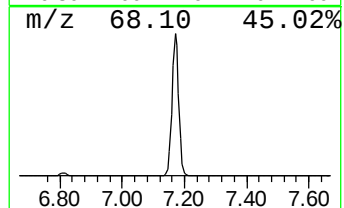
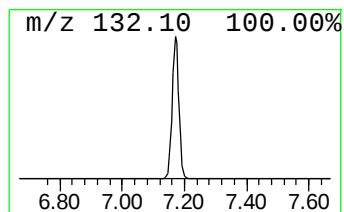
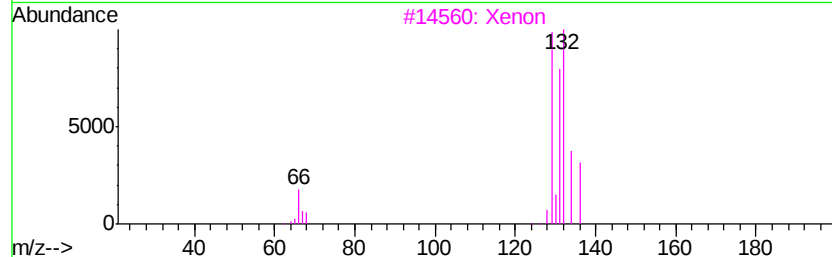
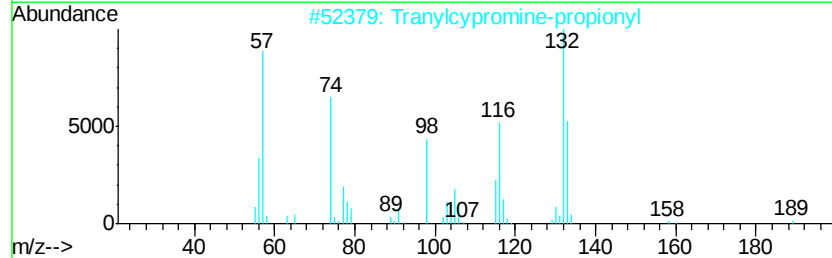
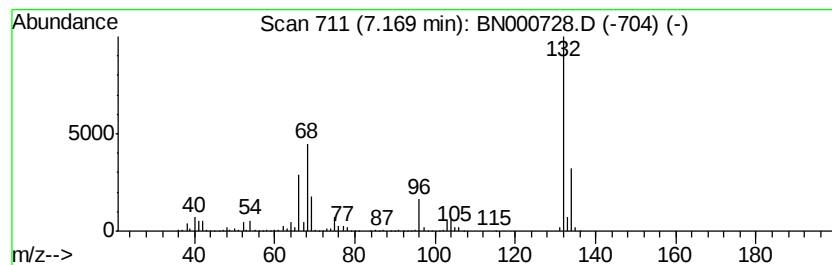
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Peak Number 2 unknown7.17 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.17	55.48 ng	2496830	1,4-Dichlorobenzene-d4	7.63

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	35
2		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
3		Xenon	132	Xe	007440-63-3	9
4		4-Chloro-2-fluoro-pyrimidine	132	C4H2ClFN2	051422-00-5	9
5		1-Methylpyrrolo[1,2-a]pyrazine	132	C8H8N2	064608-59-9	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	4.81	2.5	ng	110068	1	7.63	900025	20.0
unknown7.17	7.17	55.5	ng	2496830	1	7.63	900025	20.0