

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF041318\
 Data File : BF104491.D
 Acq On : 13 Apr 2018 20:55
 Operator : JU/SJ
 Sample : J2314-01
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 CA-AQIDW01-20180409

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:\HPCHEM1\BNA_F\METHODS\8270-BF040618.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	5.010	493	497	503	rVB	67442	72650	2.19%	0.348%
2	5.422	563	567	571	rBV	1330163	1251252	37.70%	6.001%
3	6.440	736	740	752	rBV	916687	801563	24.15%	3.844%
4	6.581	759	764	767	rBV	2641976	2371806	71.47%	11.375%
5	6.810	799	803	806	rBV	880770	679792	20.48%	3.260%
6	6.969	825	830	833	rBV	2758717	2431210	73.26%	11.660%
7	7.375	894	899	902	rBV	1780444	1892339	57.02%	9.076%
8	8.092	1017	1021	1024	rBV	1111644	882361	26.59%	4.232%
9	9.169	1199	1204	1207	rBV	3348843	3318749	100.00%	15.917%
10	9.563	1267	1271	1274	rBV	143813	119674	3.61%	0.574%
11	9.775	1304	1307	1310	rBV	54687	39103	1.18%	0.188%
12	9.851	1316	1320	1329	rVB	1049538	906168	27.30%	4.346%
13	10.275	1389	1392	1395	rVB2	65608	51876	1.56%	0.249%
14	10.639	1450	1454	1458	rBV	1575767	1556371	46.90%	7.464%
15	10.751	1470	1473	1475	rBV	57118	53703	1.62%	0.258%
16	10.775	1475	1477	1482	rVB	46117	46011	1.39%	0.221%
17	11.339	1569	1573	1576	rBV	959738	766316	23.09%	3.675%
18	12.927	1839	1843	1847	rBV	2491073	2431634	73.27%	11.662%
19	13.980	2019	2022	2026	rVB	643229	601988	18.14%	2.887%
20	15.451	2267	2272	2282	rVB	425109	575896	17.35%	2.762%

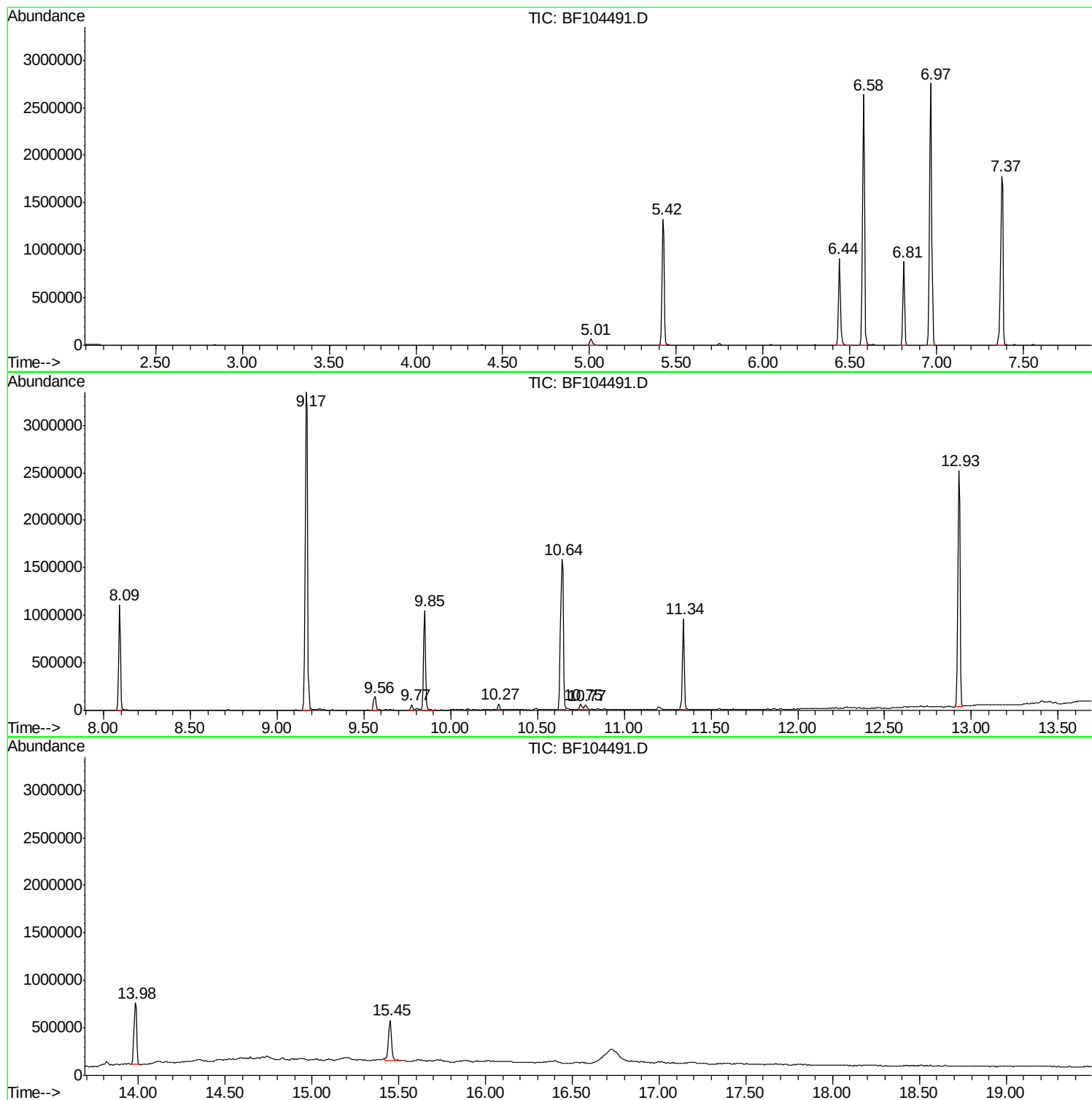
Sum of corrected areas: 20850462

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Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF040618.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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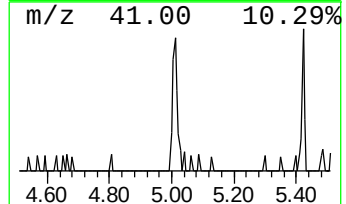
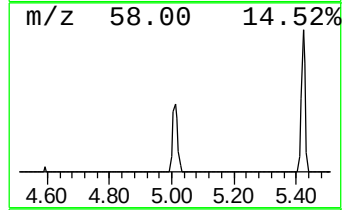
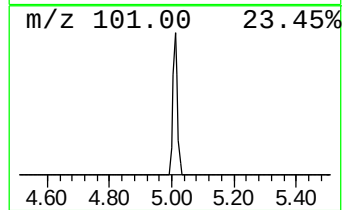
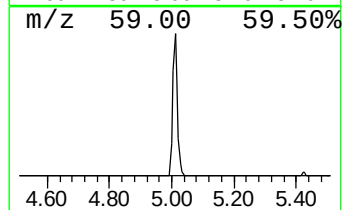
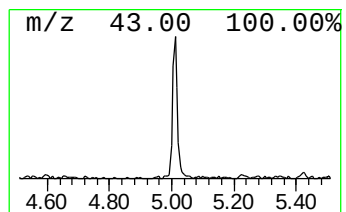
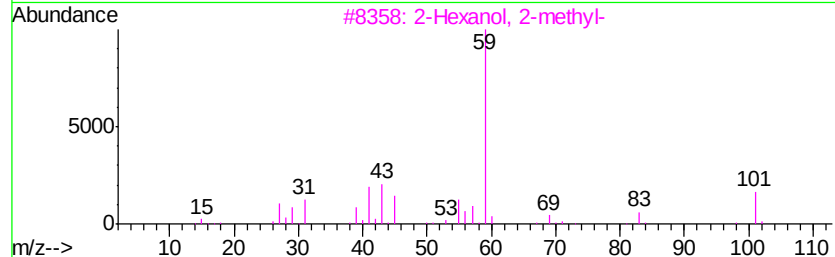
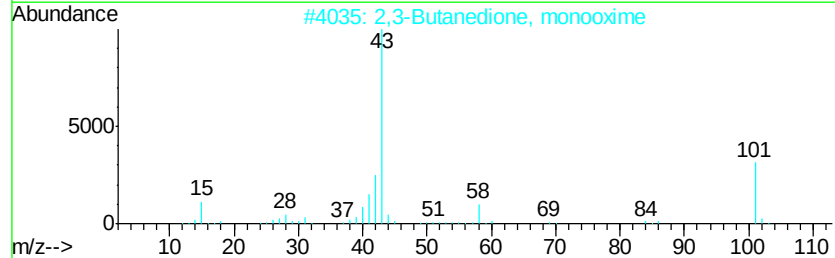
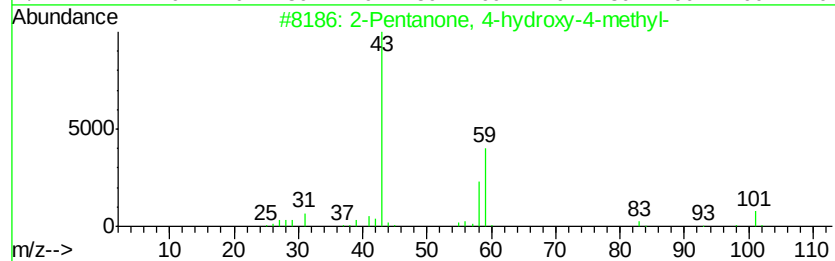
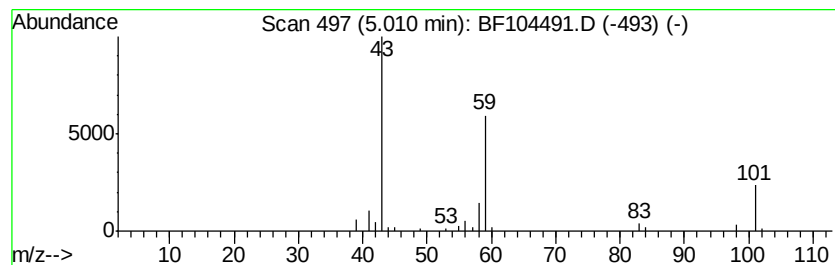
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TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.01	2.14 ng	72650	1,4-Dichlorobenzene-d4	6.81

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	59
2			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	38
3			2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	25
4			3-Pentanol	88	C5H12O	000584-02-1	23
5			Butyl aldoxime, 3-methyl-, syn-	101	C5H11NO	005780-40-5	16



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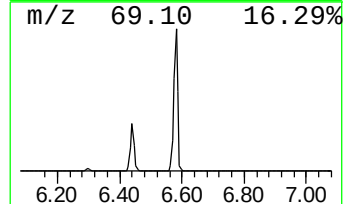
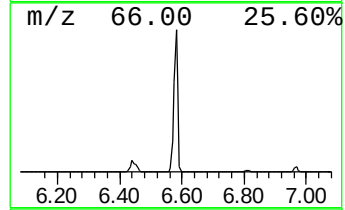
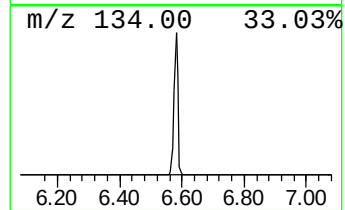
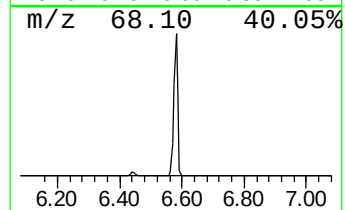
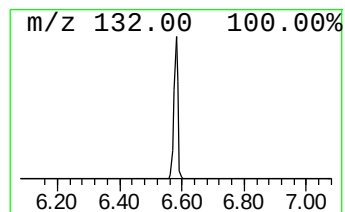
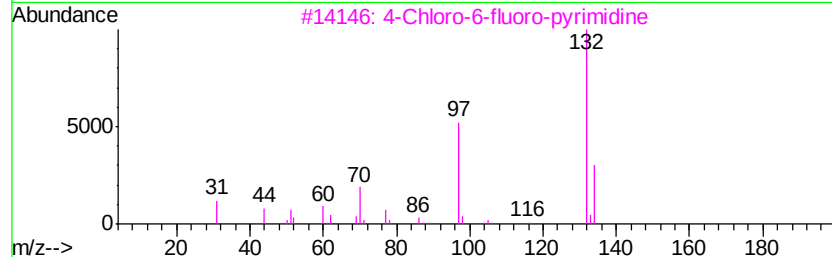
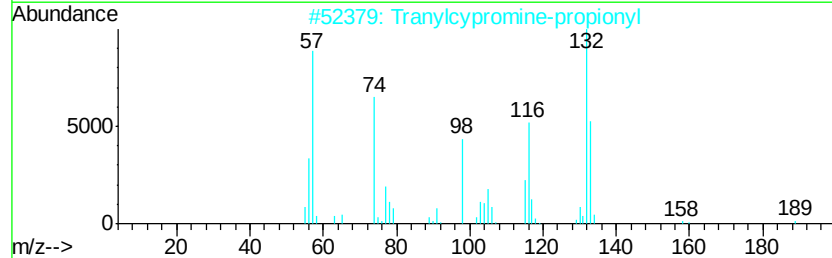
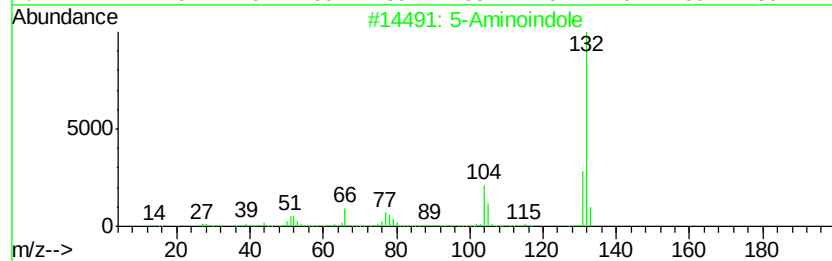
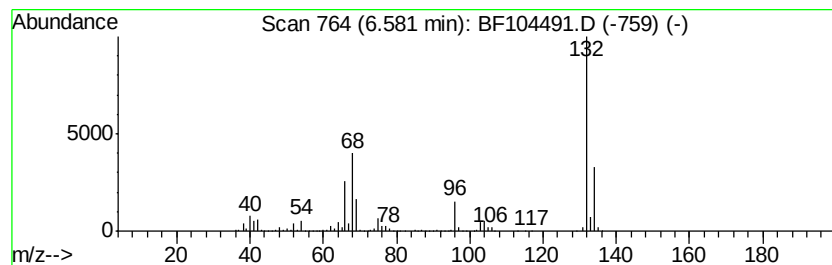
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Peak Number 2 unknown6.58 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.58	69.78 ng	2371810	1,4-Dichlorobenzene-d4	6.81

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			5-Aminoindole	132	C8H8N2	005192-03-0	12
2			Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
3			4-Chloro-6-fluoro-pyrimidine	132	C4H2ClFN2	051422-01-6	9
4			(5-Methyl-2-pyridyl)acetonitrile	132	C8H8N2	1000241-93-9	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...	5.01	2.1	ng	72650	1	6.81	679792	20.0
unknown6.58	6.58	69.8	ng	2371810	1	6.81	679792	20.0