

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA_F\DATA\BF062717\
 Data File : BF096224.D
 Acq On : 27 Jun 2017 18:57
 Operator : SJ/MA
 Sample : I3868-10 5X
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 CL-01-062217-D

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_F\Methods\8270-BF062717.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	4.963	485	489	494	rBV	174132	178303	11.63%	1.180%
2	5.381	556	560	564	rBV	1261474	1067255	69.61%	7.061%
3	6.398	729	733	736	rBV	1196142	1047974	68.35%	6.933%
4	6.539	753	757	761	rBV	1201121	1039743	67.82%	6.879%
5	6.781	794	798	801	rBV	1346691	1140769	74.41%	7.547%
6	6.934	820	824	827	rBV	942515	753663	49.16%	4.986%
7	7.333	887	892	895	rBV	658560	591064	38.55%	3.910%
8	7.416	903	906	911	rVB3	18158	19511	1.27%	0.129%
9	8.057	1011	1015	1019	rBV	1668121	1533161	100.00%	10.143%
10	9.133	1193	1198	1201	rBV	1327346	1012511	66.04%	6.699%
11	9.527	1261	1265	1267	rBV	178539	152726	9.96%	1.010%
12	9.645	1276	1285	1287	rBV6	14488	28008	1.83%	0.185%
13	9.698	1291	1294	1299	rVB2	23952	35230	2.30%	0.233%
14	9.810	1308	1313	1317	rBV2	1672483	1431269	93.35%	9.469%
15	10.222	1380	1383	1386	rBV3	17936	22987	1.50%	0.152%
16	10.257	1386	1389	1393	rVB	26513	23172	1.51%	0.153%
17	10.586	1442	1445	1449	rVB	516406	448917	29.28%	2.970%
18	10.727	1466	1469	1476	rVB3	38219	45372	2.96%	0.300%
19	11.174	1542	1545	1549	rVB3	21941	22104	1.44%	0.146%
20	11.286	1560	1564	1567	rBV	1222867	1106384	72.16%	7.320%
21	11.310	1567	1568	1571	rVB	122524	73614	4.80%	0.487%
22	11.363	1571	1577	1579	rBV	27133	26170	1.71%	0.173%
23	11.521	1600	1604	1608	rVB3	24245	28195	1.84%	0.187%
24	11.786	1647	1649	1651	rBV2	18129	16377	1.07%	0.108%
25	11.821	1651	1655	1658	rVV3	54403	56503	3.69%	0.374%
26	11.898	1665	1668	1671	rBV	33521	34750	2.27%	0.230%
27	12.080	1695	1699	1701	rBV2	19335	22106	1.44%	0.146%
28	12.492	1766	1769	1773	rBV	200995	179296	11.69%	1.186%
29	12.604	1786	1788	1793	rBV5	46010	55059	3.59%	0.364%
30	12.721	1802	1808	1814	rBV	178789	187284	12.22%	1.239%
31	12.868	1830	1833	1837	rBV	715778	611008	39.85%	4.042%
32	13.051	1858	1864	1868	rBV2	26407	36492	2.38%	0.241%
33	13.115	1873	1875	1878	rVB3	20036	19435	1.27%	0.129%
34	13.274	1900	1902	1906	rBV4	16443	17564	1.15%	0.116%

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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-BF062717.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

35	13.774	1985	1987	1992	rVB4	44424	39666	2.59%	0.262%
36	13.915	2006	2011	2014	rBV	1187138	1104369	72.03%	7.306%
37	13.939	2014	2015	2018	rVB	73316	47958	3.13%	0.317%
38	15.339	2248	2253	2259	rBV	706948	859260	56.04%	5.685%

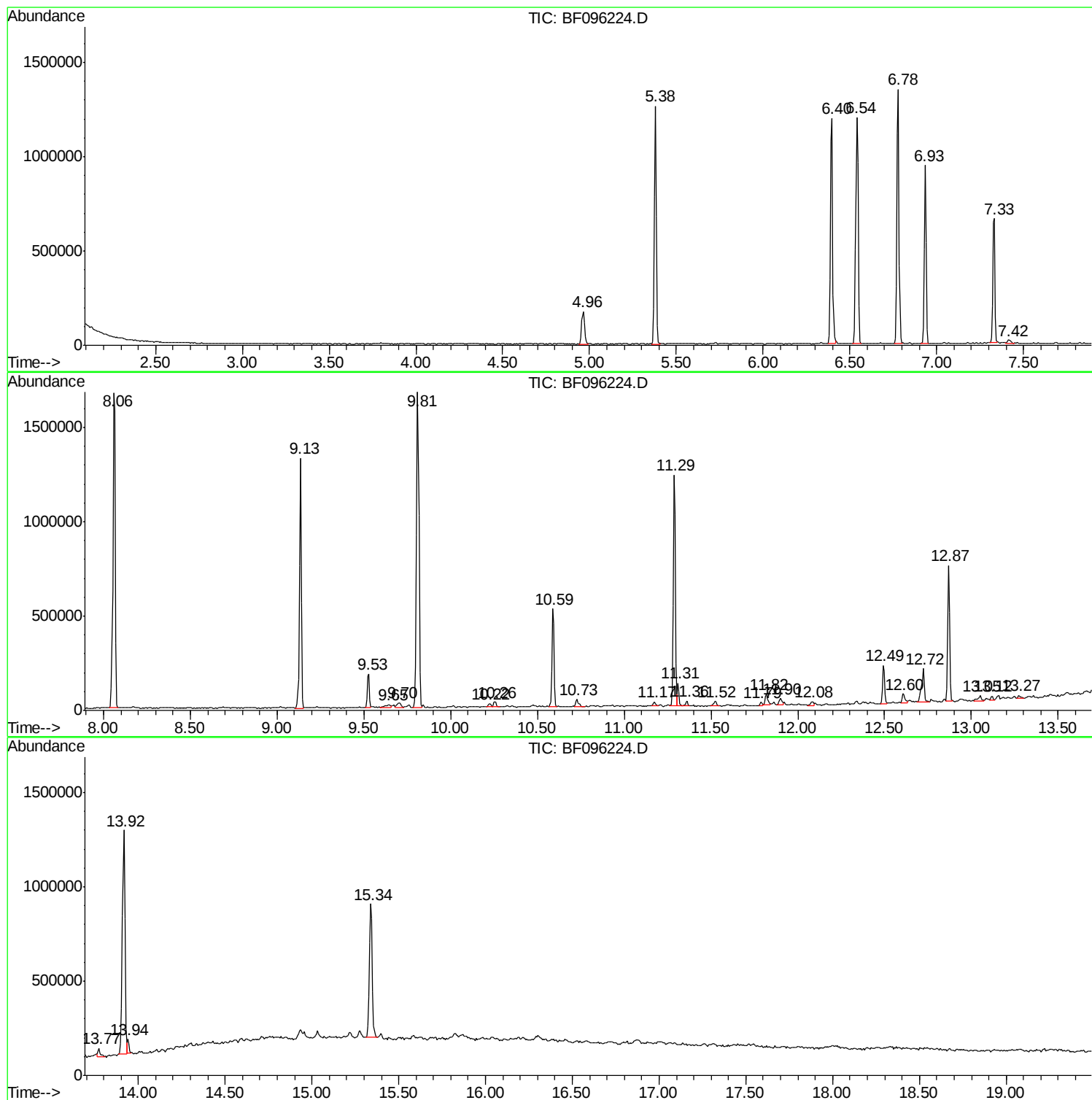
Sum of corrected areas: 15115229

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Quant Method : Z:\HPCHEM1\BNA_F\Methods\8270-BF062717.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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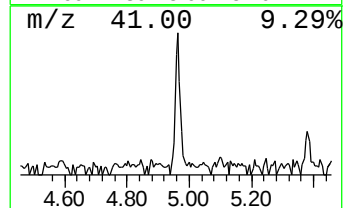
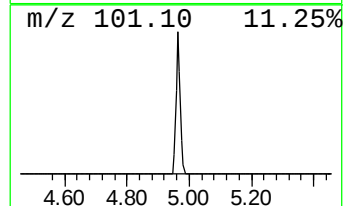
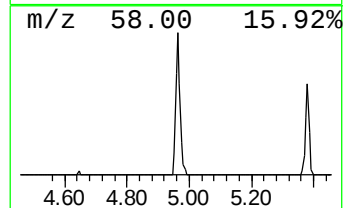
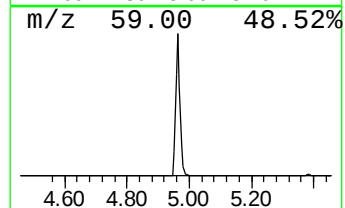
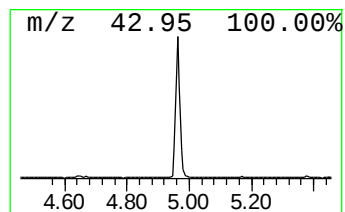
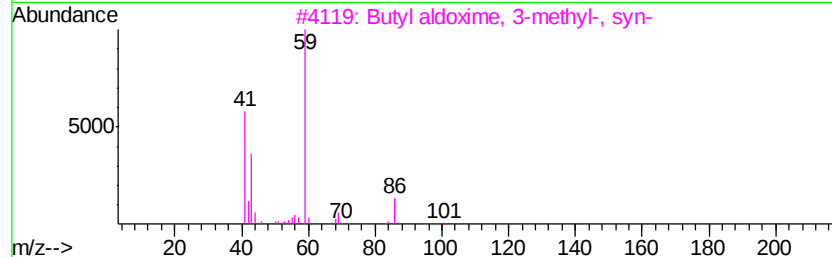
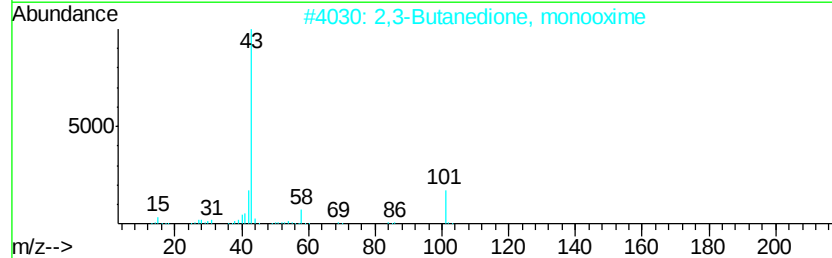
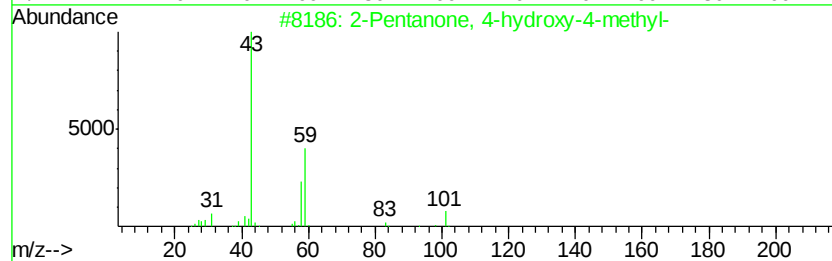
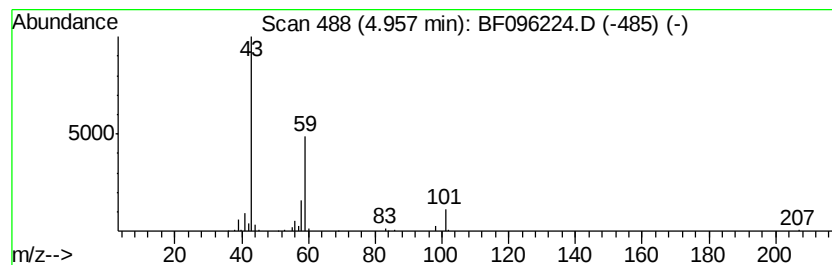
Quant Method : Z:\HPCHEM1\BNA_F\Methods\8270-BF062717.M
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.96	3.13 ng	178303	1,4-Dichlorobenzene-d4	6.78

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50
2			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	27
3			Butyl aldoxime, 3-methyl-, syn-	101	C5H11NO	005780-40-5	23
4			1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	10
5			Butane	58	C4H10	000106-97-8	9



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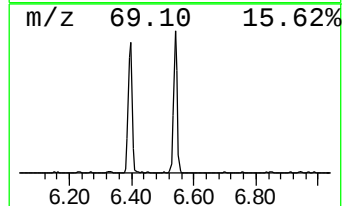
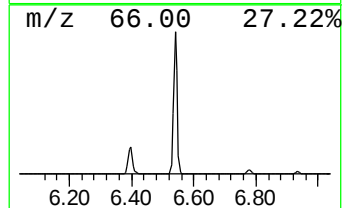
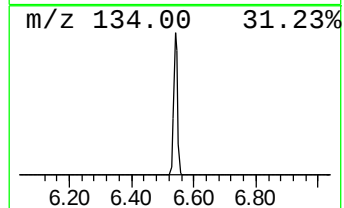
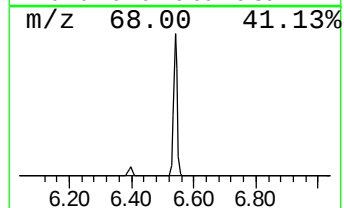
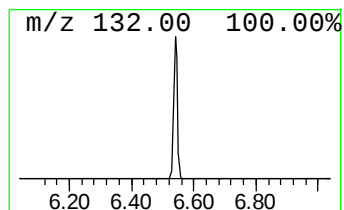
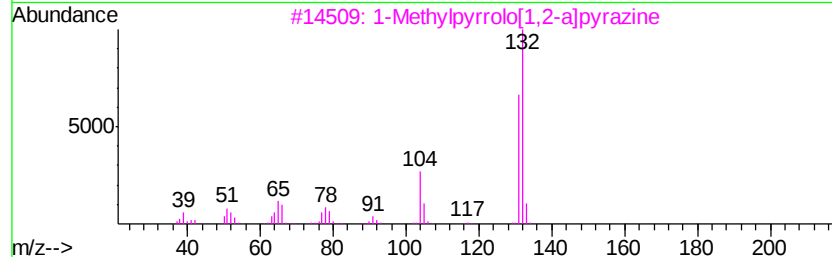
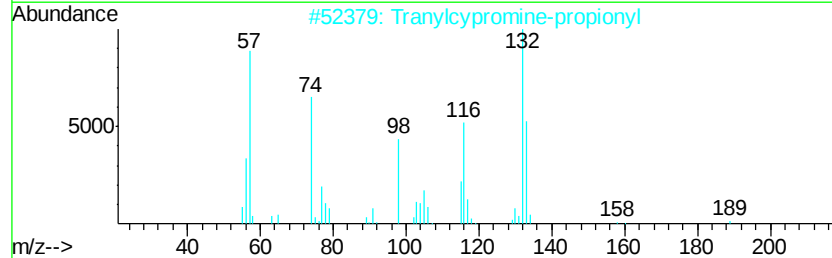
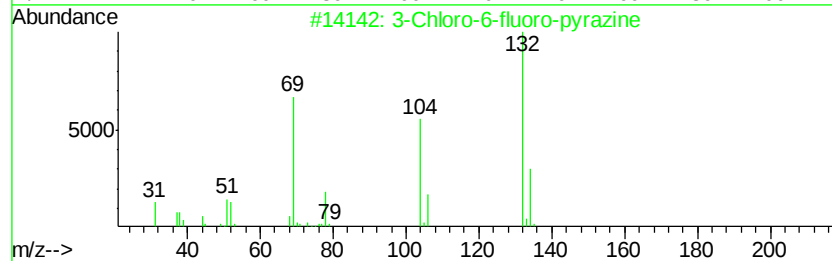
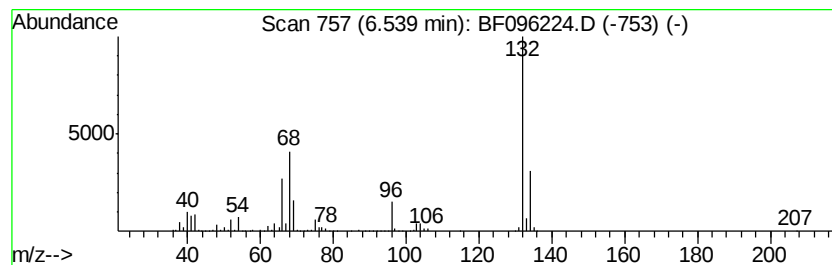
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Peak Number 2 unknown6.54 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.54	18.23 ng	1039740	1,4-Dichlorobenzene-d4	6.78

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2			Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
3			1-Methylpyrrolo[1,2-a]pyrazine	132	C8H8N2	064608-59-9	9
4			5-Aminoindole	132	C8H8N2	005192-03-0	9
5			4-Methylpyrrolo[1,2-a]pyrazine	132	C8H8N2	064608-60-2	9



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TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	4.96	3.1	ng	178303	1	6.78	1140770	20.0
unknown6.54	6.54	18.2	ng	1039740	1	6.78	1140770	20.0