

Data Path : Z:\HPCHEM1\BNA F\DATA\BF062817\
 Data File : BF096256.D
 Acq On : 28 Jun 2017 18:24
 Operator : SJ/MA
 Sample : I3852-09 5X
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 G-27-0-1.5

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	496	rBV	232717	229356	16.81%	1.503%
2	5.387	557	561	564	rBV	1257666	1062970	77.89%	6.966%
3	6.404	730	734	739	rBV	1234824	1100405	80.63%	7.211%
4	6.545	754	758	762	rBV	1325916	1077439	78.95%	7.061%
5	6.781	795	798	802	rVB	1273027	991211	72.63%	6.496%
6	6.940	821	825	828	rVB	940888	802354	58.79%	5.258%
7	7.334	888	892	895	rBV	855638	633276	46.40%	4.150%
8	7.745	959	962	966	rBV2	22607	21666	1.59%	0.142%
9	8.063	1012	1016	1019	rBV	1712055	1364752	100.00%	8.944%
10	8.304	1054	1057	1061	rVB	42221	31131	2.28%	0.204%
11	9.139	1195	1199	1202	rVB	1244589	1085679	79.55%	7.115%
12	9.234	1212	1215	1218	rBV3	19278	21624	1.58%	0.142%
13	9.528	1262	1265	1268	rBV	142299	109736	8.04%	0.719%
14	9.757	1302	1304	1310	rBV	29477	24496	1.79%	0.161%
15	9.816	1310	1314	1317	rBV	1589922	1325828	97.15%	8.688%
16	10.257	1386	1389	1392	rVB2	39509	29325	2.15%	0.192%
17	10.592	1443	1446	1450	rBV	560120	473989	34.73%	3.106%
18	10.728	1466	1469	1476	rBV2	47254	66576	4.88%	0.436%
19	11.175	1537	1545	1549	rBV5	39346	69833	5.12%	0.458%
20	11.210	1549	1551	1556	rVB3	24876	23730	1.74%	0.156%
21	11.263	1556	1560	1562	rBV3	55807	55974	4.10%	0.367%
22	11.292	1562	1565	1568	rVV2	1339196	1167517	85.55%	7.651%
23	11.316	1568	1569	1572	rVV	114043	65692	4.81%	0.430%
24	11.363	1572	1577	1580	rVV4	31883	34726	2.54%	0.228%
25	11.604	1615	1618	1620	rBV	22337	20226	1.48%	0.133%
26	11.822	1652	1655	1659	rBV2	70299	73116	5.36%	0.479%
27	11.904	1666	1669	1671	rBV3	33036	35811	2.62%	0.235%
28	12.010	1685	1687	1690	rBV2	19939	19317	1.42%	0.127%
29	12.345	1741	1744	1748	rBV5	35019	34948	2.56%	0.229%
30	12.498	1767	1770	1773	rBV	144316	136130	9.97%	0.892%
31	12.610	1787	1789	1792	rBV3	29573	29987	2.20%	0.197%
32	12.727	1806	1809	1812	rVB	147092	120353	8.82%	0.789%
33	12.874	1831	1834	1838	rVB	992581	807927	59.20%	5.295%
34	13.739	1979	1981	1983	rBV	49249	33322	2.44%	0.218%

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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	rVB2	83695	85350	6.25%	0.559%
36	13.921	2008	2012	2015	rBV	1178841	1080367	79.16%	7.080%
37	14.104	2041	2043	2046	rBV	67747	54742	4.01%	0.359%
38	15.351	2251	2255	2260	rVB	741201	858709	62.92%	5.627%

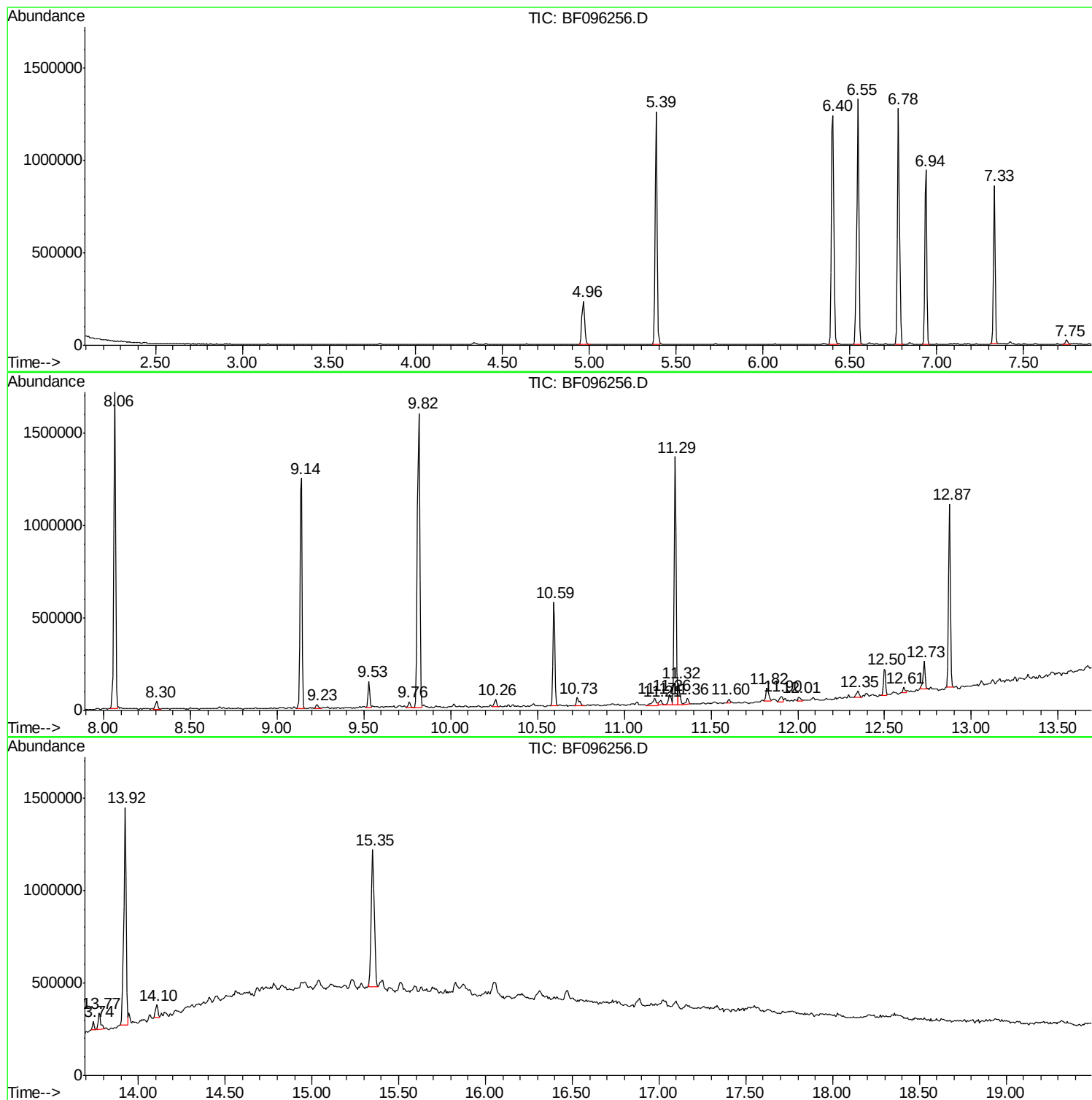
Sum of corrected areas: 15259590

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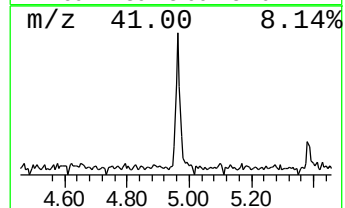
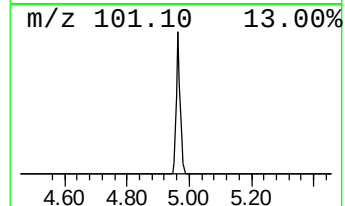
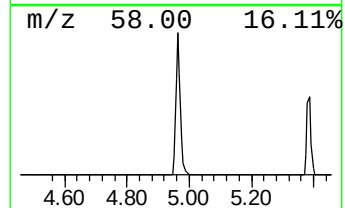
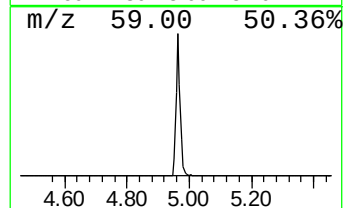
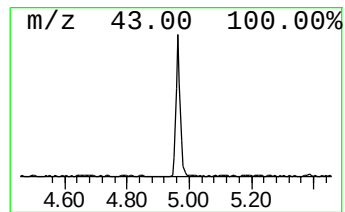
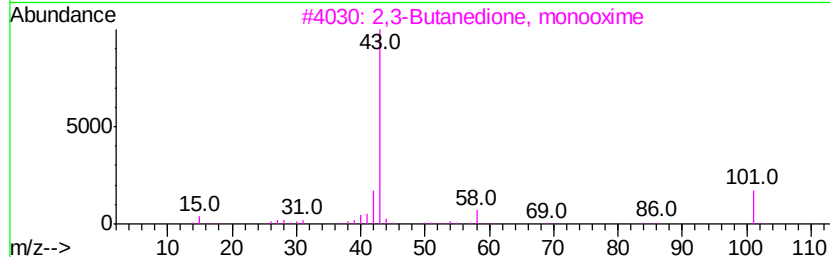
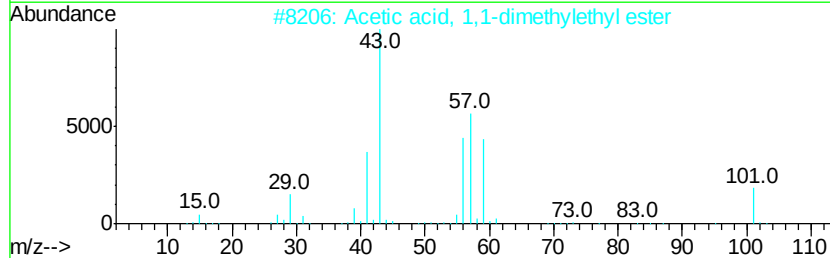
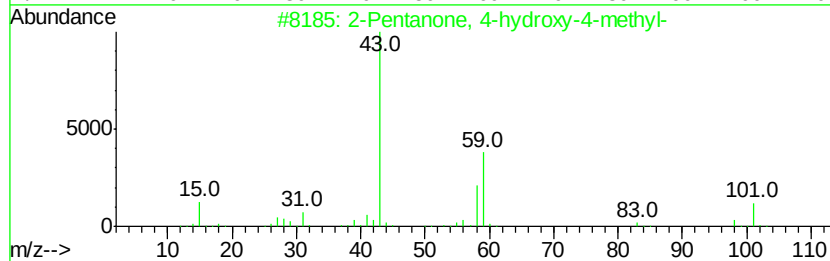
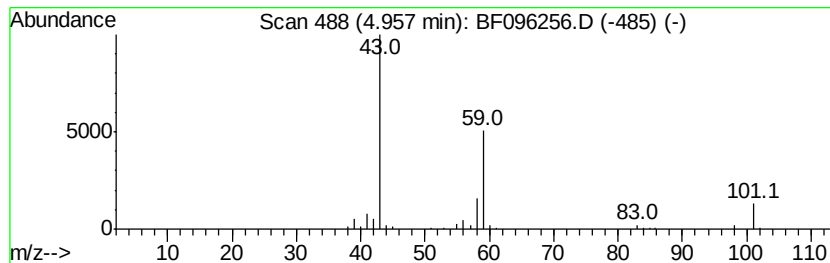
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.96	4.63 ng	229356	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			Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	38
3			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9
4			1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	9
5			1,2-Butanediol	90	C4H10O2	000584-03-2	9



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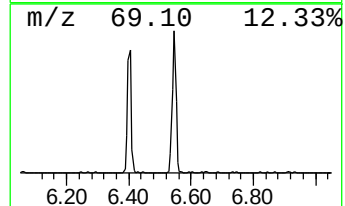
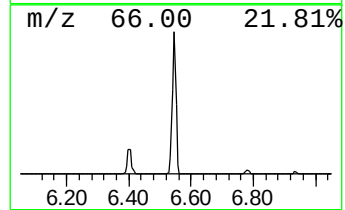
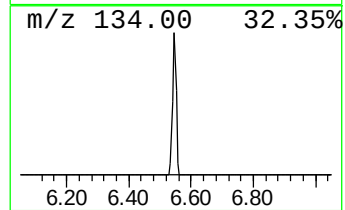
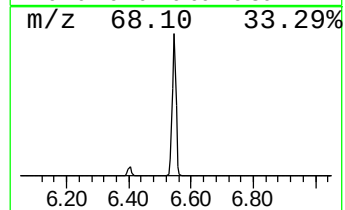
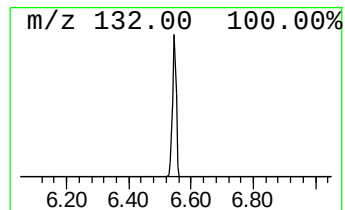
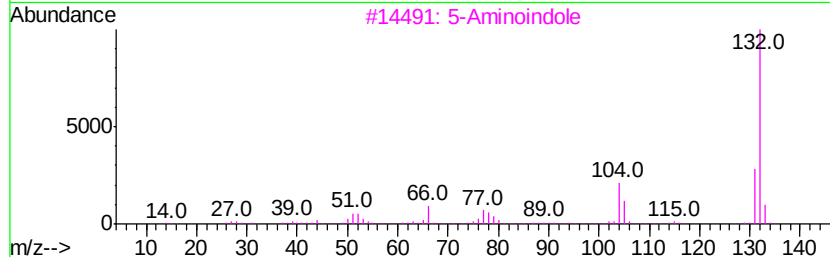
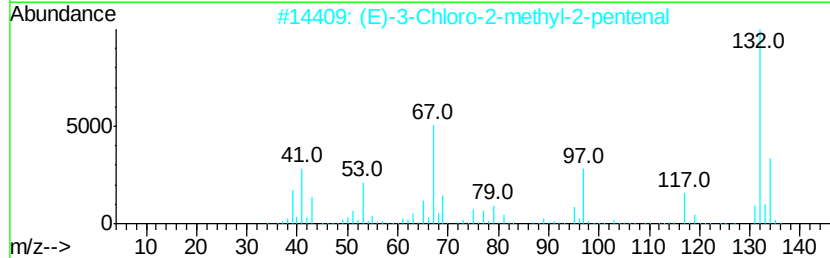
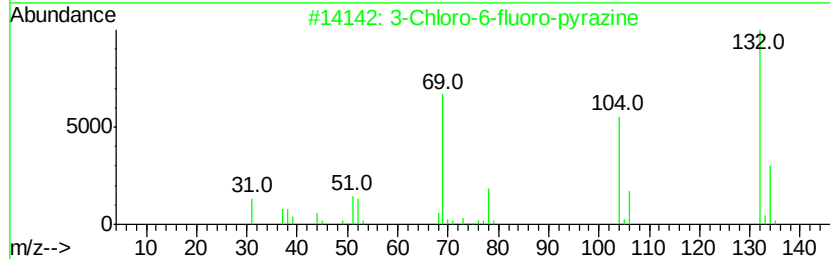
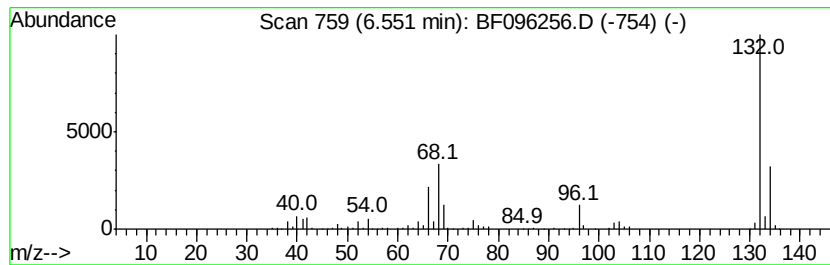
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Peak Number 2 unknown6.55 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.55	21.74 ng	1077440	1,4-Dichlorobenzene-d4	6.78

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	16
3		5-Aminoindole	132	C8H8N2	005192-03-0	12
4		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
5		Amphetaminil	250	C17H18N2	017590-01-1	10



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	4.96	4.6	ng	229356	1	6.78	991211	20.0
unknown6.55	6.55	21.7	ng	1077440	1	6.78	991211	20.0