

Data Path : Z:\HPCHEM1\BNA F\DATA\BF031318\
 Data File : BF103602.D
 Acq On : 13 Mar 2018 14:21
 Operator : SJ/JU
 Sample : PB107288BL
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

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-BF022218.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC

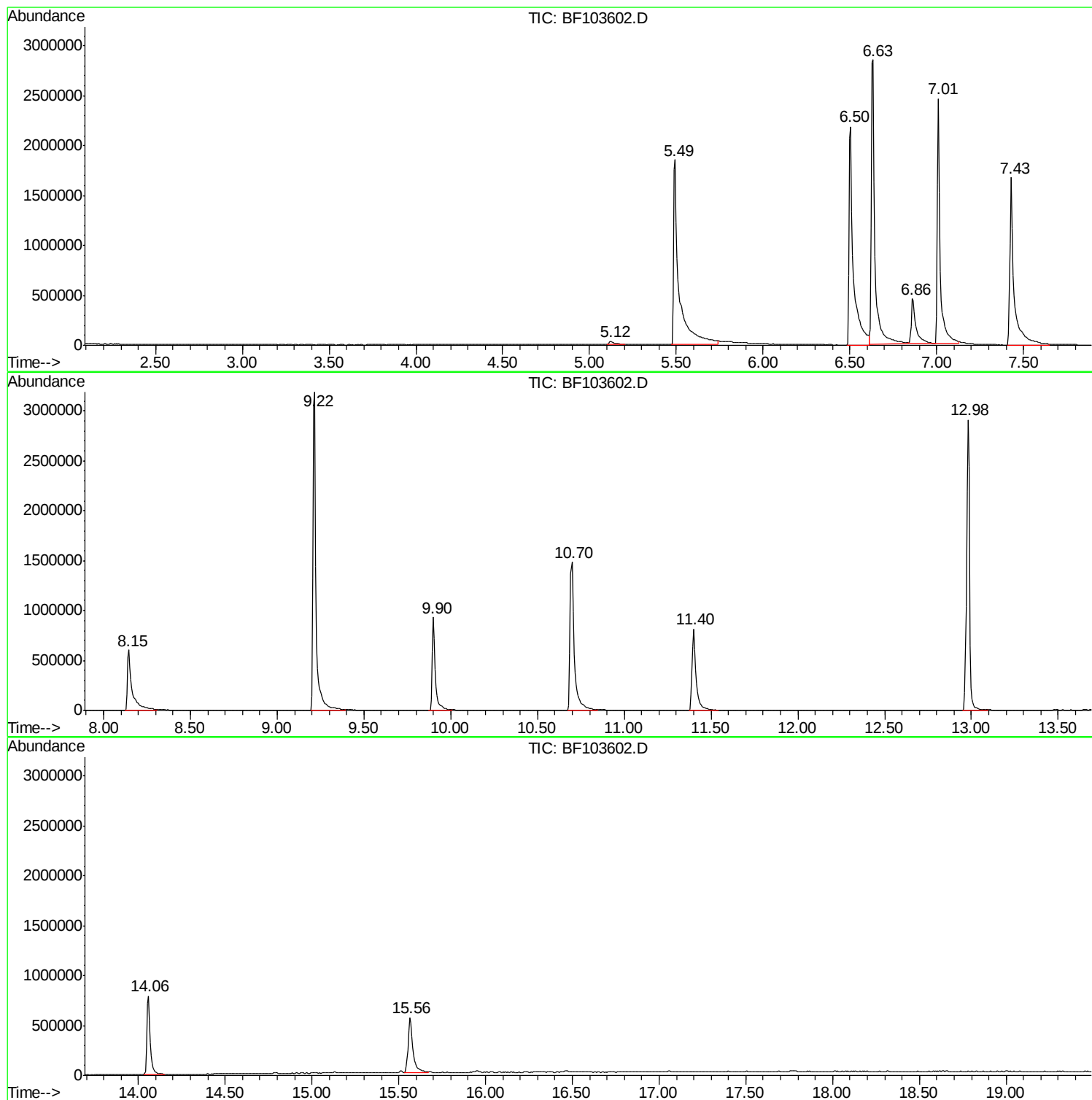
peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.122	512	516	530	rBV2	33490	92151	2.13%	0.261%
2	5.492	575	579	621	rBV	1853452	4242614	98.15%	12.038%
3	6.504	747	751	769	rBV	2186907	3960968	91.63%	11.239%
4	6.633	769	773	808	rVV	2846704	4322777	100.00%	12.265%
5	6.857	808	811	833	rVB	448951	846881	19.59%	2.403%
6	7.010	833	837	857	rBV	2455184	3012701	69.69%	8.548%
7	7.428	905	908	946	rBV	1679688	2829010	65.44%	8.027%
8	8.145	1026	1030	1057	rBV	604505	1186981	27.46%	3.368%
9	9.216	1208	1212	1242	rBV	3188009	4149229	95.99%	11.773%
10	9.898	1324	1328	1346	rBV2	929455	1246054	28.83%	3.535%
11	10.698	1460	1464	1489	rBV	1482891	2355668	54.49%	6.684%
12	11.398	1579	1583	1607	rBV	813683	1272078	29.43%	3.609%
13	12.980	1847	1852	1871	rBV	2906471	3512762	81.26%	9.967%
14	14.057	2031	2035	2051	rBV	791033	1131956	26.19%	3.212%
15	15.562	2286	2291	2310	rBV	551733	1082393	25.04%	3.071%

Sum of corrected areas: 35244223

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Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF022218.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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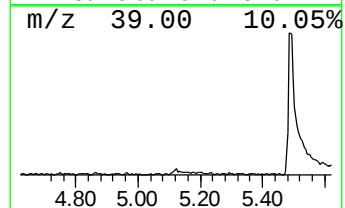
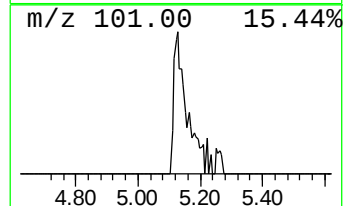
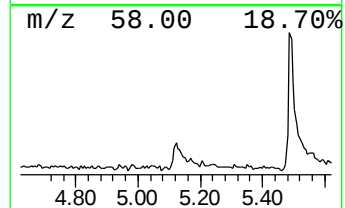
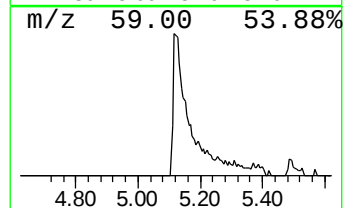
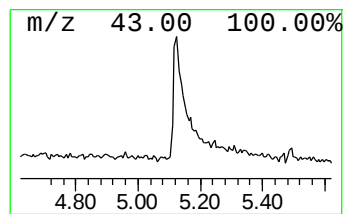
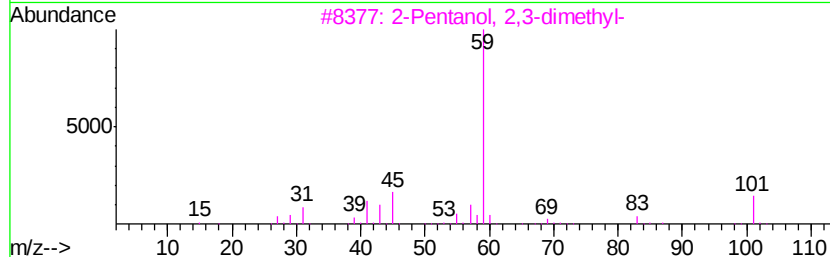
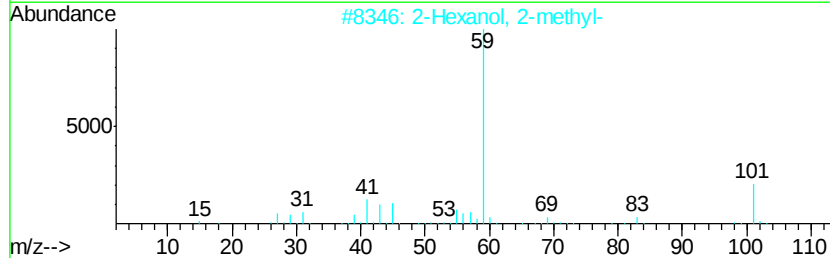
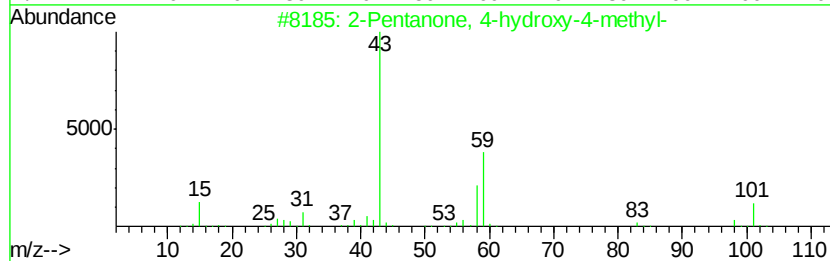
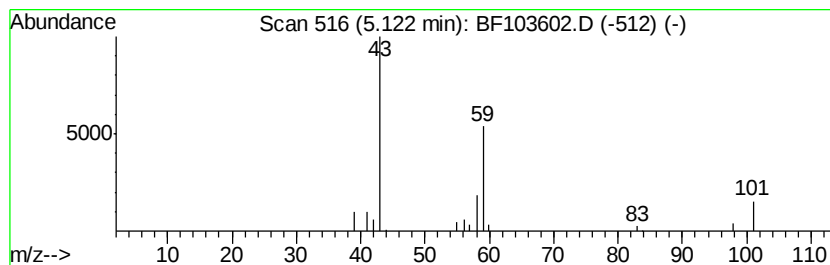
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.12	2.18 ng	92151	1,4-Dichlorobenzene-d4	6.86

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	64
2			2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	50
3			2-Pentanol, 2,3-dimethyl-	116	C7H16O	004911-70-0	9
4			3-Octanol	130	C8H18O	000589-98-0	9
5			2-Pentanol, 2,4-dimethyl-	116	C7H16O	000625-06-9	9



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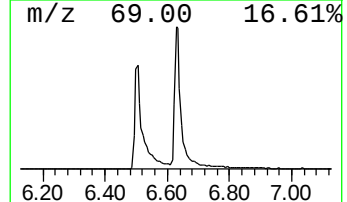
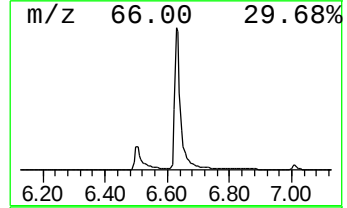
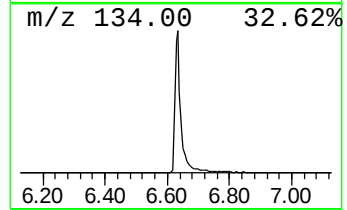
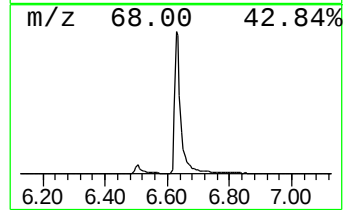
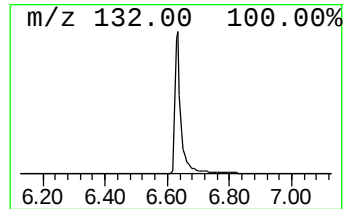
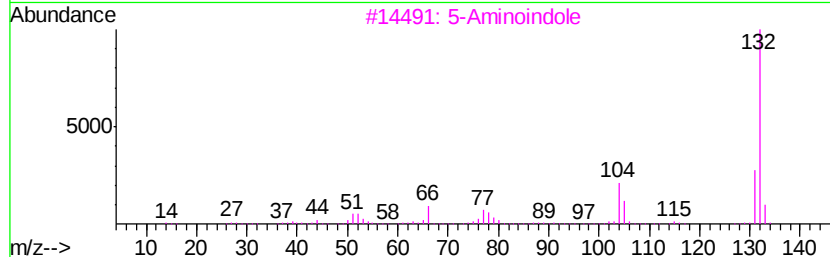
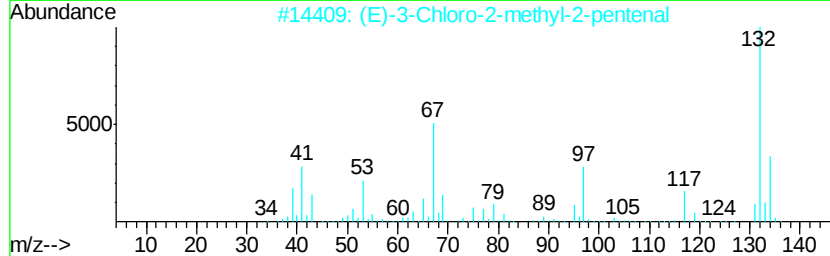
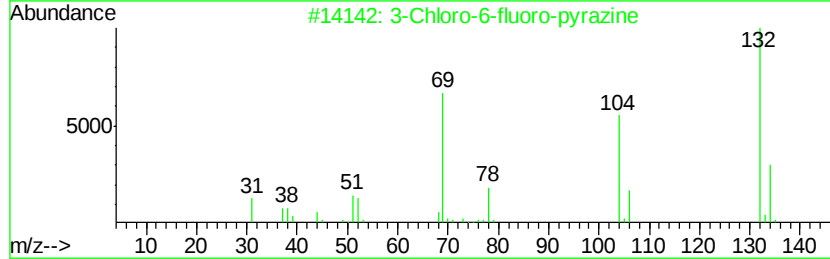
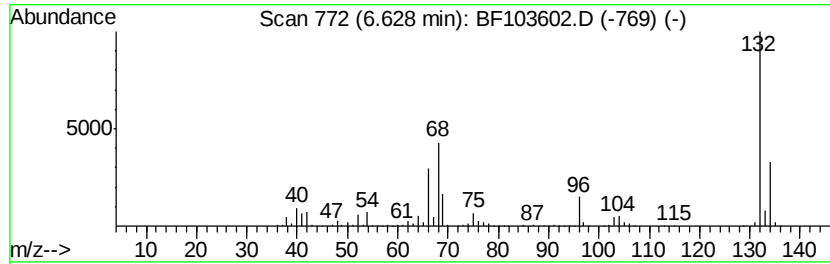
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Peak Number 2 unknown6.63 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.63	102.09 ng	4322780	1,4-Dichlorobenzene-d4	6.86

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	17
3		5-Aminoindole	132	C8H8N2	005192-03-0	12
4		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
5		Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.12	2.2	ng	92151	1	6.86	846881	20.0
unknown6.63	6.63	102.1	ng	4322780	1	6.86	846881	20.0