

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF010218\
 Data File : BF101844.D
 Acq On : 2 Jan 2018 12:08
 Operator : SJ/JU
 Sample : PB105401BL
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB105401BL

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-BF121417.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	5.004	493	496	516	rBV	269906	522309	2.04%	0.904%
2	5.404	561	564	593	rBV	2714976	3796154	14.82%	6.571%
3	6.198	688	699	734	rBV3	5200858	25614413	100.00%	44.339%
4	6.428	734	738	755	rVV	2767606	3781749	14.76%	6.546%
5	6.551	755	759	769	rVB	3141500	3458910	13.50%	5.987%
6	6.775	794	797	812	rVB	732667	757795	2.96%	1.312%
7	6.934	820	824	842	rVB	2660317	2802969	10.94%	4.852%
8	7.351	891	895	915	rBV	2172750	2461603	9.61%	4.261%
9	8.057	1012	1015	1029	rBV	952584	1017589	3.97%	1.761%
10	9.134	1193	1198	1201	rBV	3919802	3981891	15.55%	6.893%
11	9.816	1310	1314	1321	rVB	1127392	1092232	4.26%	1.891%
12	10.616	1446	1450	1463	rBV	2183774	2301981	8.99%	3.985%
13	11.310	1564	1568	1577	rBV	1200372	1148139	4.48%	1.987%
14	12.892	1832	1837	1840	rBV	3551871	3601963	14.06%	6.235%
15	13.963	2015	2019	2029	rVB	755986	794283	3.10%	1.375%
16	15.433	2264	2269	2277	rBV	422871	636047	2.48%	1.101%

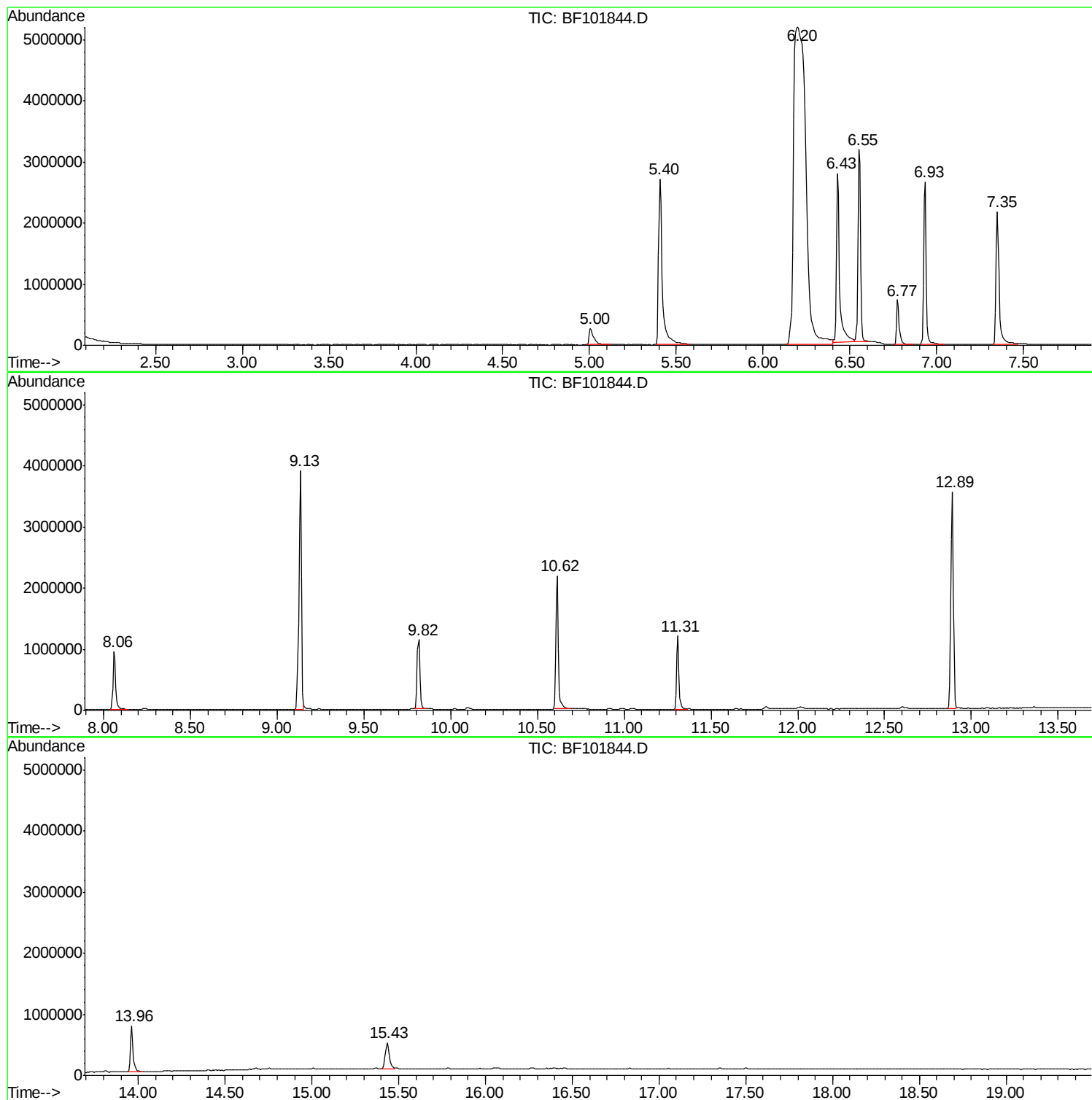
Sum of corrected areas: 57770027

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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF121417.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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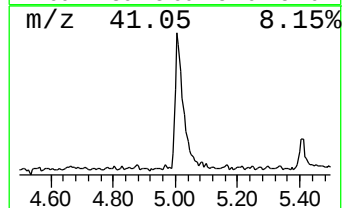
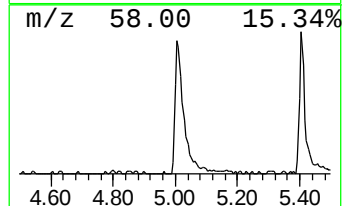
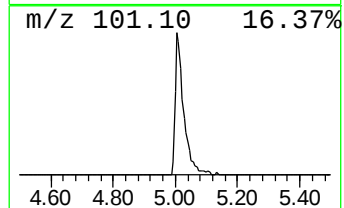
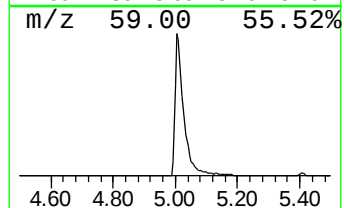
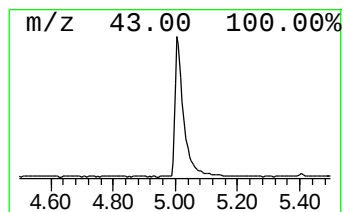
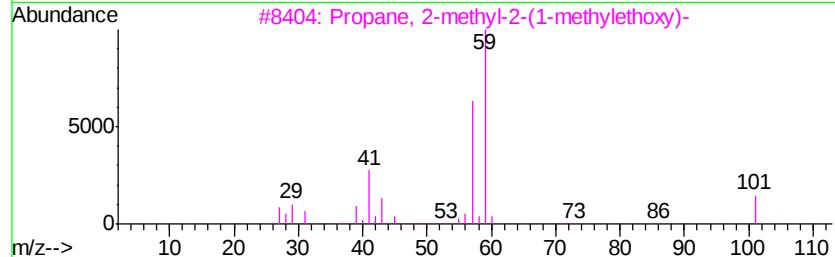
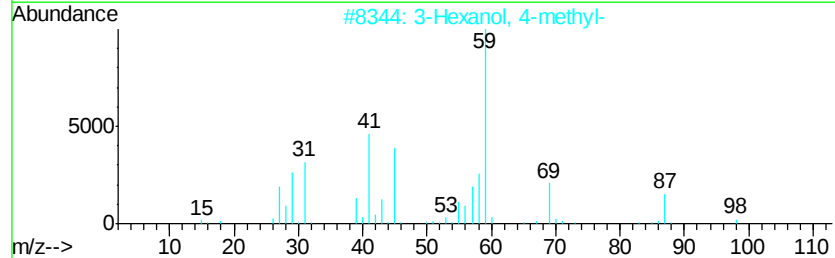
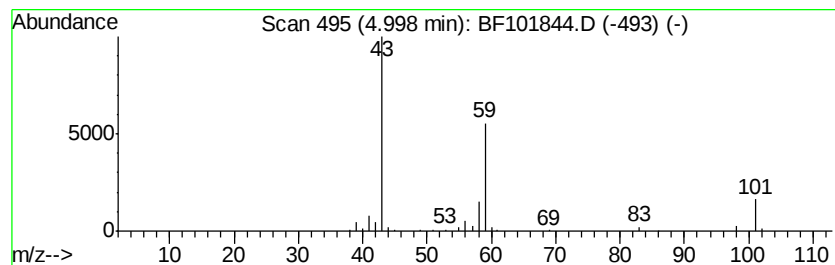
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.00	13.79 ng	522309	1,4-Dichlorobenzene-d4	6.77

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	45
2		3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	33
3		Propane, 2-methyl-2-(1-methyleth...	116	C7H16O	017348-59-3	33
4		Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	32
5		1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	10



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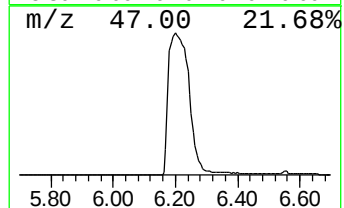
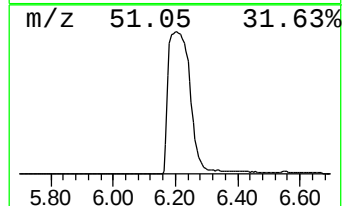
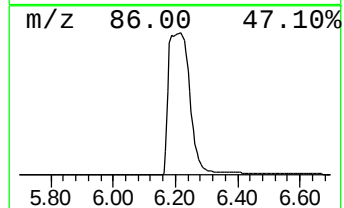
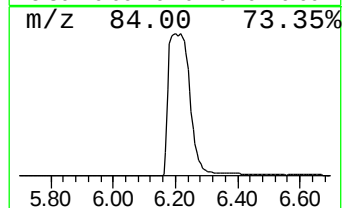
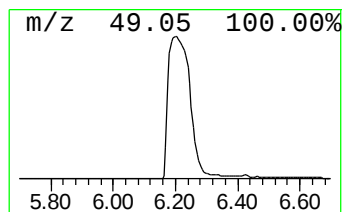
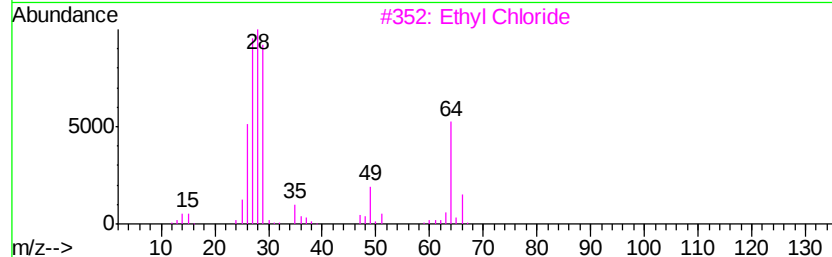
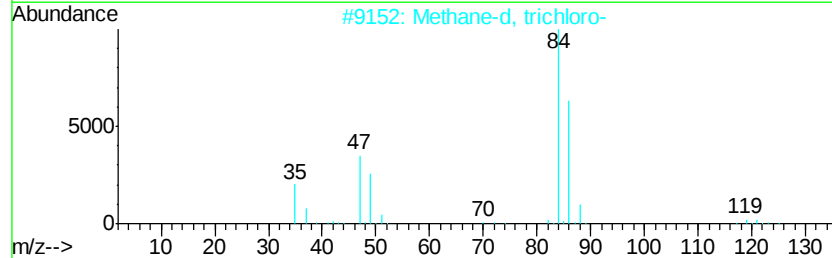
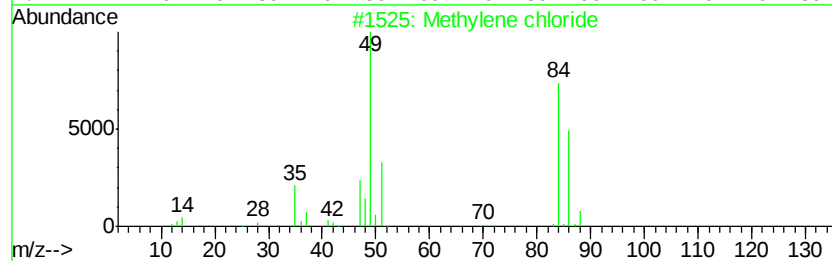
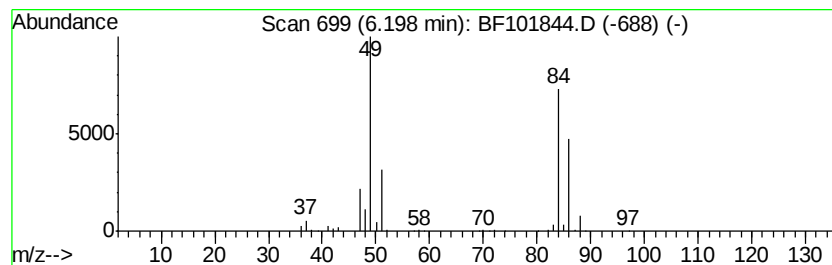
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Peak Number 2 Methylene chloride Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.20	676.03 ng	25614400	1,4-Dichlorobenzene-d4	6.77

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Methylene chloride	84	CH2Cl2	000075-09-2	95
2		Methane-d, trichloro-	119	CDCl3	000865-49-6	14
3		Ethyl Chloride	64	C2H5Cl	000075-00-3	4
4		Krypton	84	Kr	007439-90-9	3
5		Acetic acid, chloro-, ethyl ester	122	C4H7ClO2	000105-39-5	2



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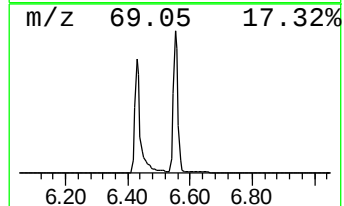
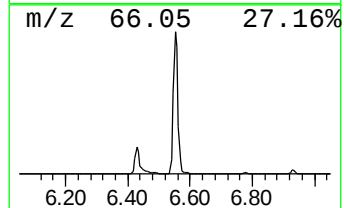
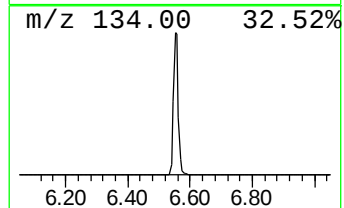
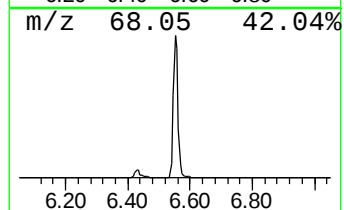
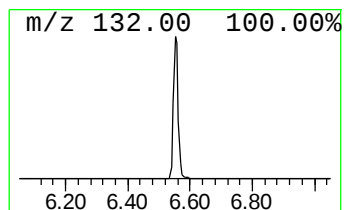
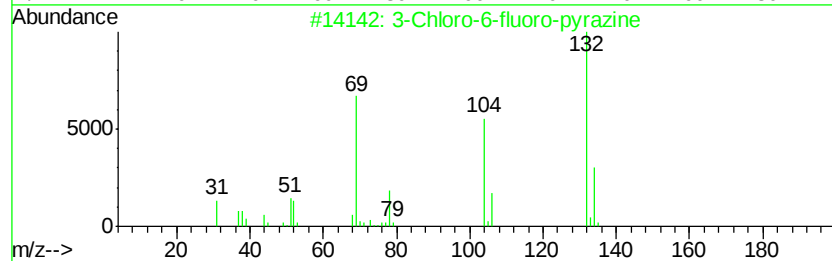
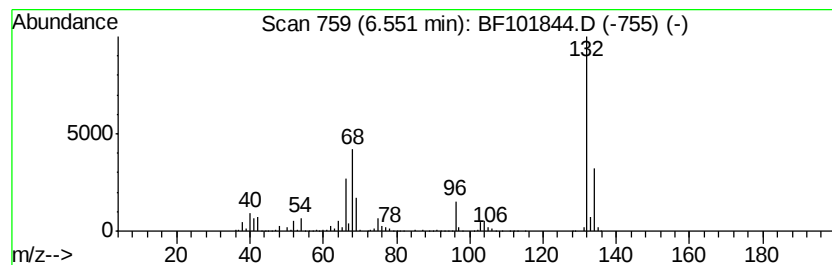
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Peak Number 3 unknown6.55 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.55	91.29 ng	3458910	1,4-Dichlorobenzene-d4	6.77

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	35
2			5-Aminoindole	132	C8H8N2	005192-03-0	12
3			Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
4			4-Chloro-2-fluoro-pyrimidine	132	C4H2ClFN2	051422-00-5	9
5			2H-Cyclopenta[d]pyridazine, 2-me...	132	C8H8N2	022291-85-6	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.00	13.8	ng	522309	1	6.77	757795	20.0
Methylene chloride	6.20	676.0	ng	25614400	1	6.77	757795	20.0
unknown6.55	6.55	91.3	ng	3458910	1	6.77	757795	20.0