

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF060117\
 Data File : BF095605.D
 Acq On : 1 Jun 2017 20:44
 Operator : SJ/MA
 Sample : I3384-03
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 MW-3

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-BF050217.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	4.775	521	525	547	rBV	242305	471737	18.61%	3.341%
2	5.486	643	646	666	rBV	128564	391533	15.44%	2.773%
3	7.122	920	924	933	rBV	91698	217185	8.57%	1.538%
4	7.198	933	937	953	rVB	557524	1083650	42.74%	7.674%
5	7.539	991	995	1006	rBV	300056	412560	16.27%	2.922%
6	7.775	1030	1035	1057	rBV	1069503	1534678	60.53%	10.868%
7	8.451	1145	1150	1185	rBV	754002	1229253	48.49%	8.705%
8	9.569	1336	1340	1362	rBV	379726	605603	23.89%	4.289%
9	11.345	1635	1642	1663	rBV	1857769	2535215	100.00%	17.953%
10	12.392	1816	1820	1838	rBV2	449442	661874	26.11%	4.687%
11	13.692	2036	2041	2061	rBV	433420	722684	28.51%	5.118%
12	14.798	2224	2229	2248	rBV	448134	668805	26.38%	4.736%
13	17.150	2623	2629	2641	rBV	2054859	2524021	99.56%	17.874%
14	18.497	2853	2858	2866	rBV	428036	617098	24.34%	4.370%
15	20.168	3138	3142	3156	rBV	295632	445456	17.57%	3.154%

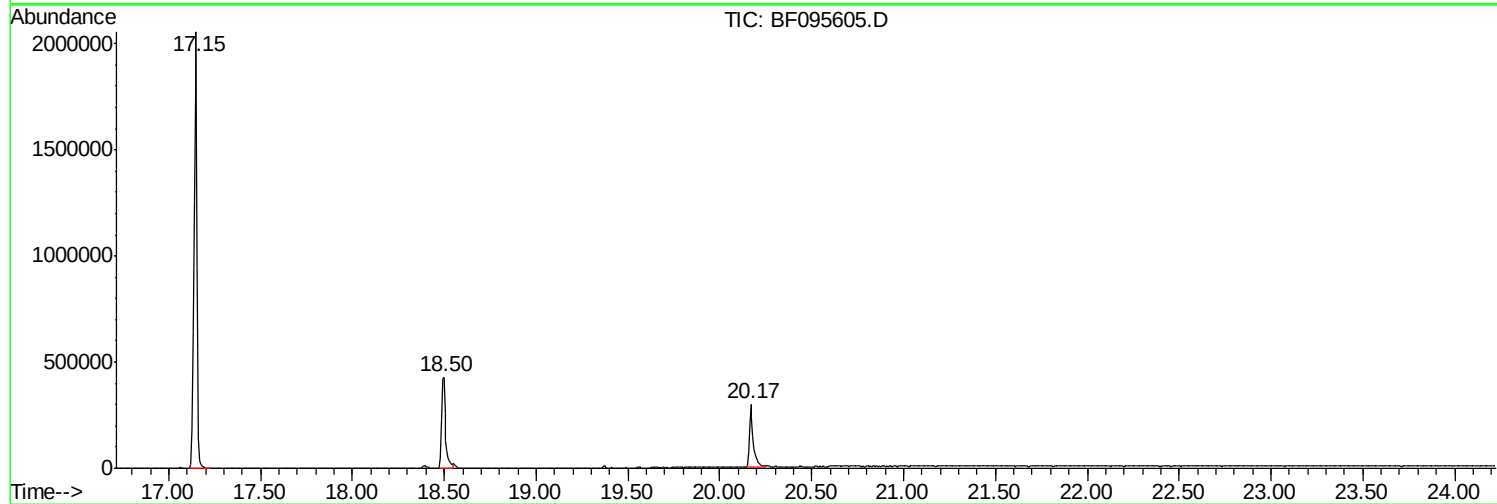
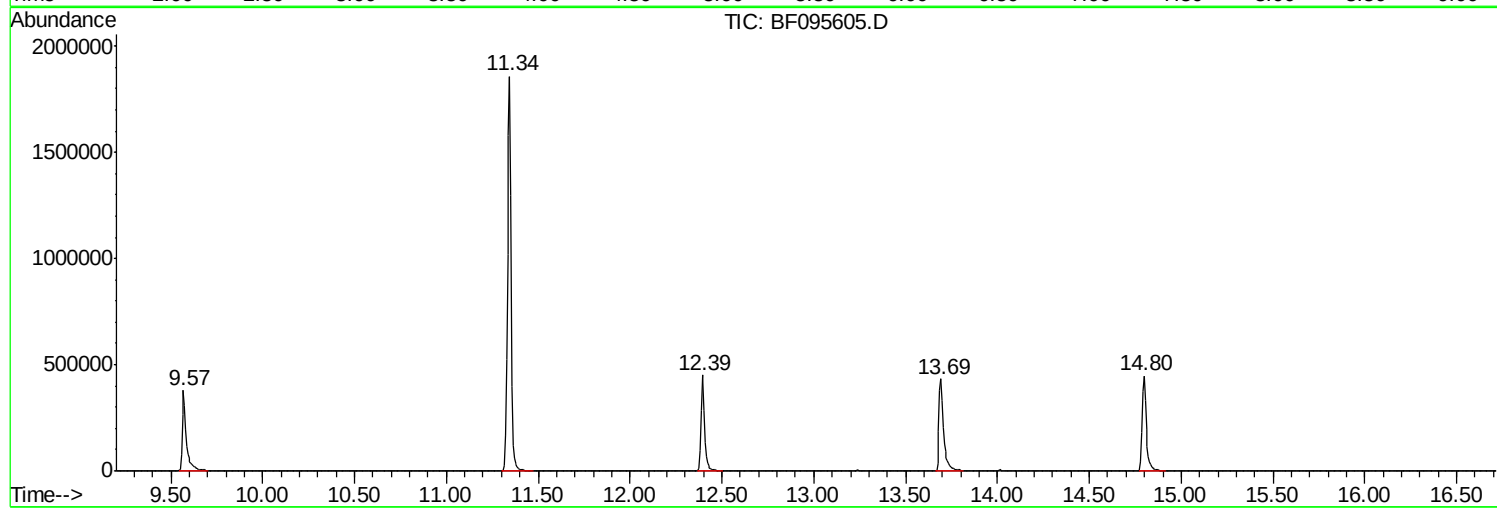
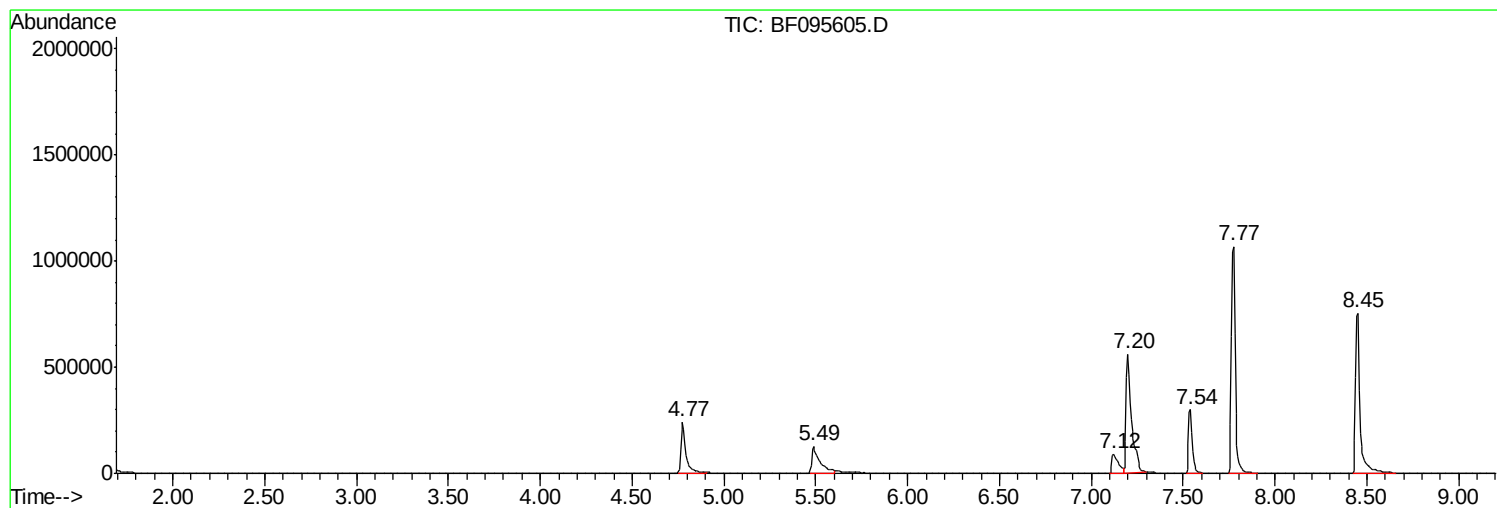
Sum of corrected areas: 14121352

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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF050217.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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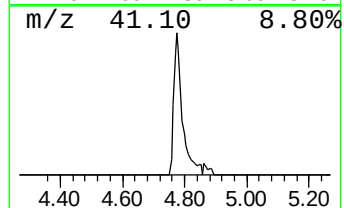
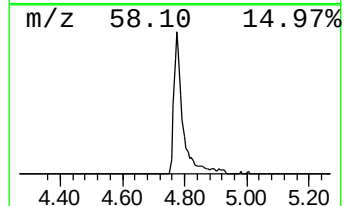
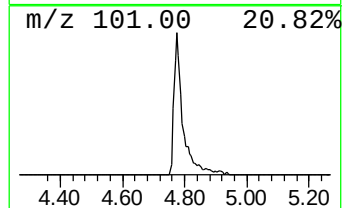
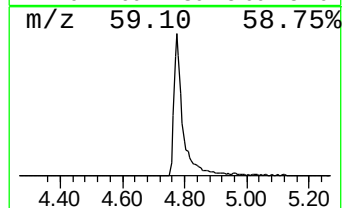
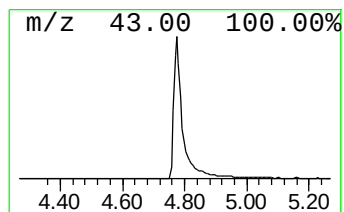
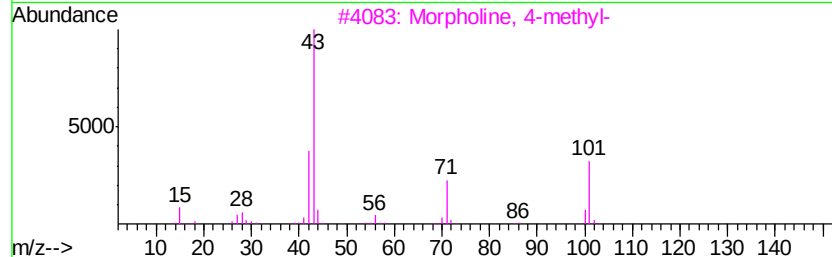
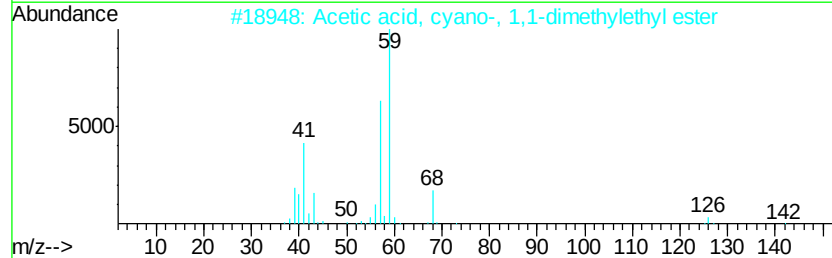
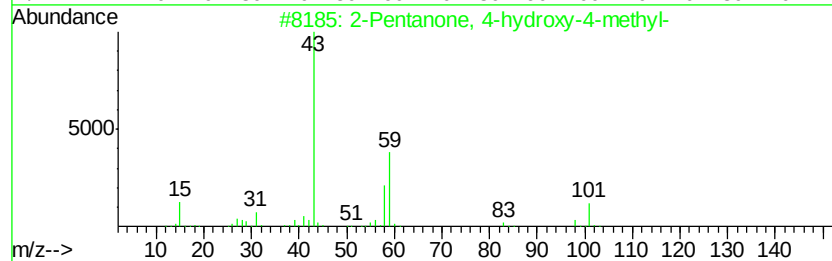
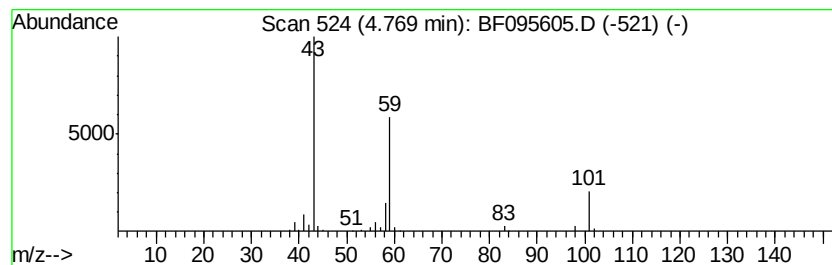
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.77	22.87 ng	471737	1,4-Dichlorobenzene-d4	7.54

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	53
2		Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	17
3		Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9
4		2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9
5		5-Hexen-2-one	98	C6H10O	000109-49-9	9



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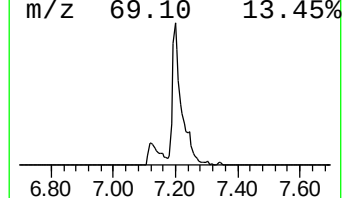
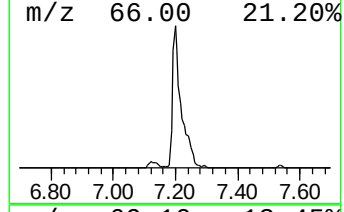
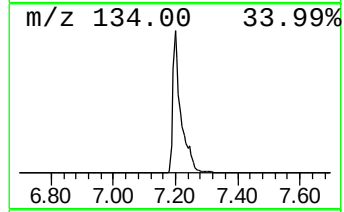
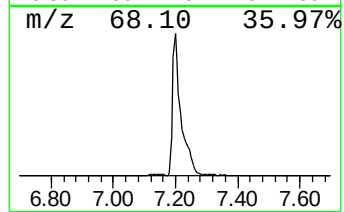
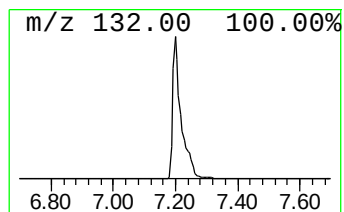
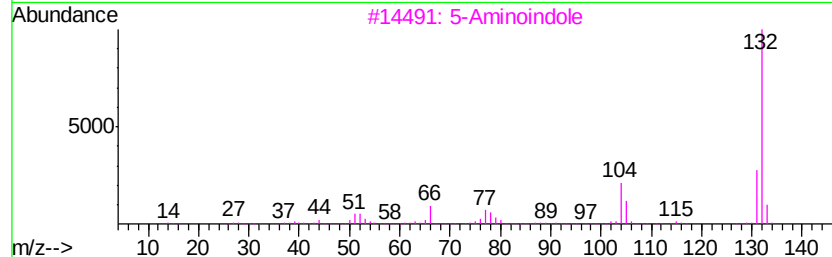
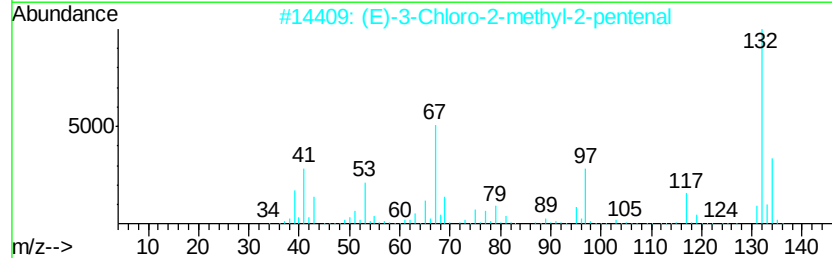
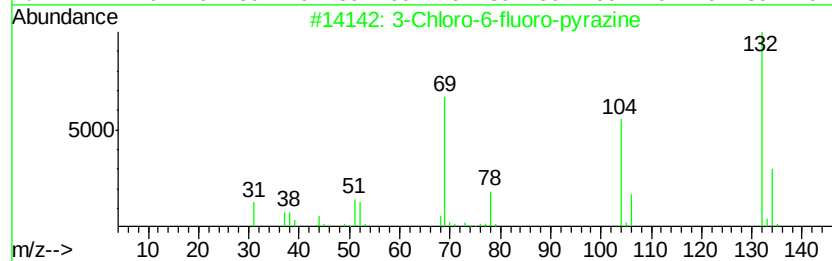
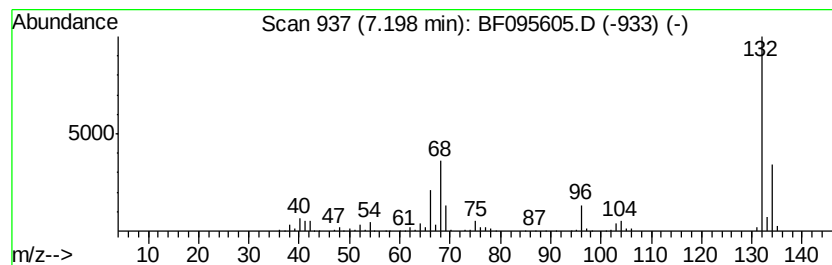
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Peak Number 2 unknown7.20 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.20	52.53 ng	1083650	1,4-Dichlorobenzene-d4	7.54

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	38
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	17
3		5-Aminoindole	132	C8H8N2	005192-03-0	12
4		1H-Indazole, 7-methyl-	132	C8H8N2	1000316-00-7	10
5		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	4.77	22.9	ng	471737	1	7.54	412560	20.0
unknown7.20	7.20	52.5	ng	1083650	1	7.54	412560	20.0