

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF102517\
 Data File : BF099822.D
 Acq On : 25 Oct 2017 19:23
 Operator : SJ/JU
 Sample : PB103580BL
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB103580BL

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-BF102317.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	5.010	494	497	514	rBV	144892	212096	11.65%	1.536%
2	5.416	563	566	587	rBV	1387033	1409180	77.43%	10.208%
3	6.434	735	739	758	rBV	1416375	1373525	75.47%	9.950%
4	6.563	758	761	773	rVB	1674342	1491129	81.93%	10.802%
5	6.793	797	800	813	rBB	456684	381390	20.96%	2.763%
6	6.945	823	826	844	rBB	1310940	1211328	66.56%	8.775%
7	7.363	893	897	917	rBV	940775	865335	47.55%	6.269%
8	8.075	1015	1018	1036	rBV	643599	529278	29.08%	3.834%
9	9.157	1196	1202	1205	rBV	1931398	1820020	100.00%	13.184%
10	9.834	1314	1317	1325	rBB	624958	579774	31.86%	4.200%
11	10.633	1449	1453	1468	rBB	991566	874178	48.03%	6.333%
12	11.328	1568	1571	1579	rBV	619014	563837	30.98%	4.084%
13	12.916	1837	1841	1844	rBV	1787250	1616442	88.81%	11.710%
14	13.980	2019	2022	2029	rBV	477176	465005	25.55%	3.369%
15	15.451	2266	2272	2279	rVB	322233	411931	22.63%	2.984%

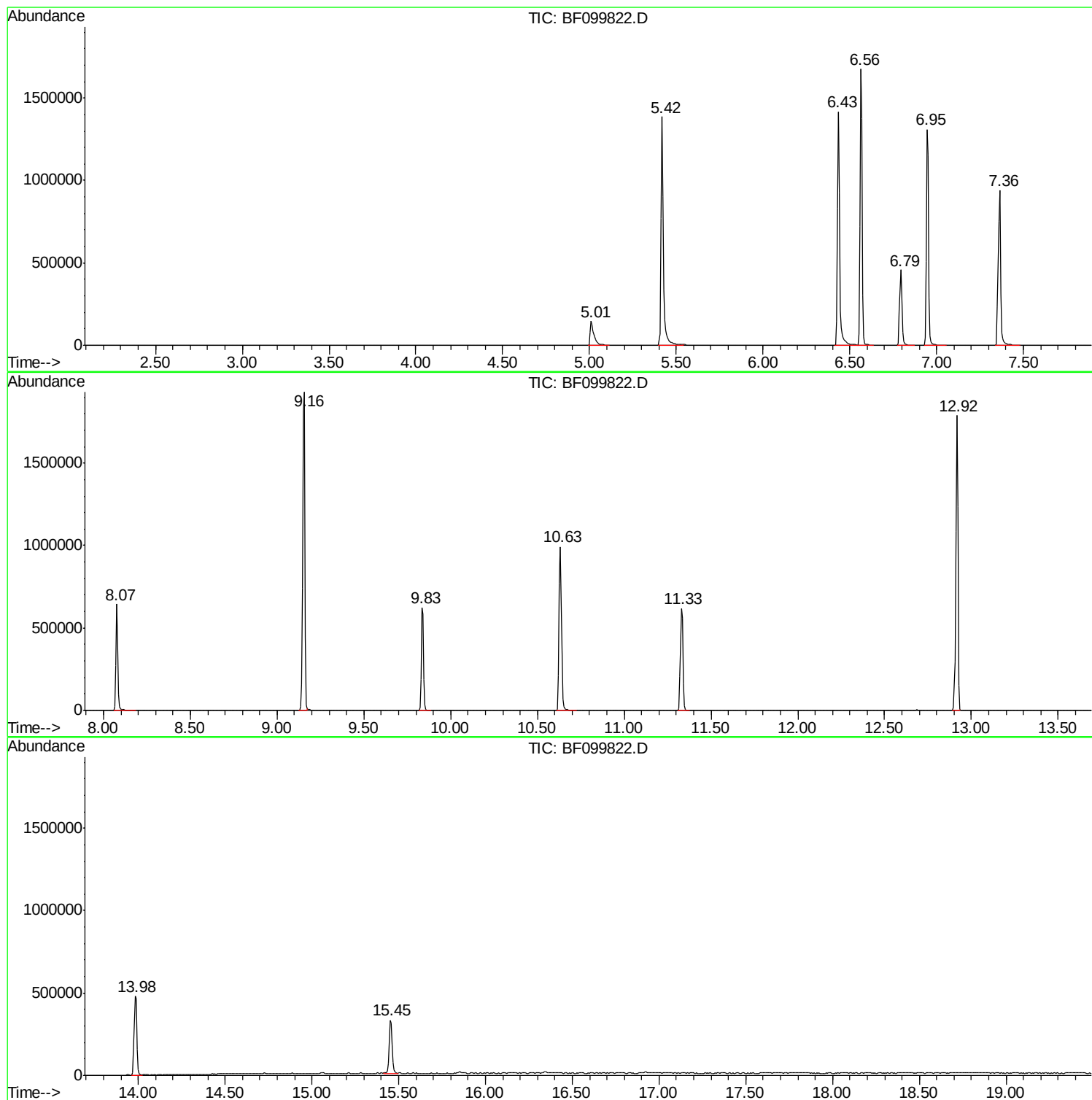
Sum of corrected areas: 13804448

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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF102317.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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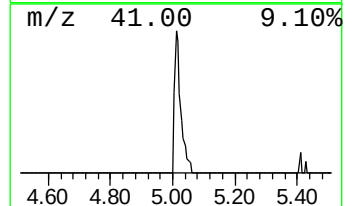
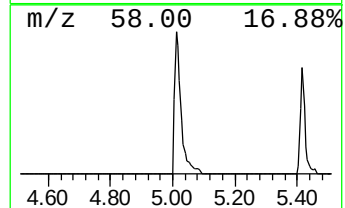
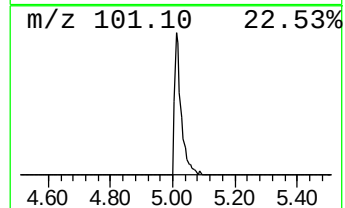
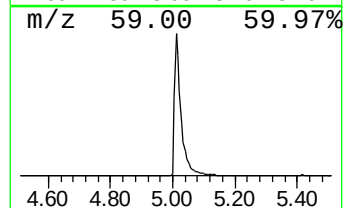
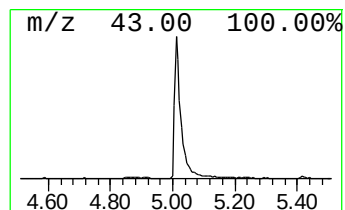
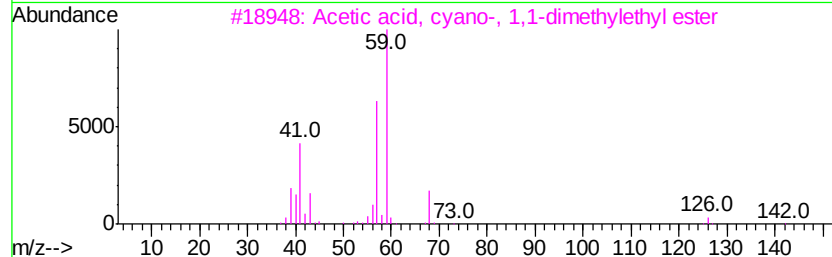
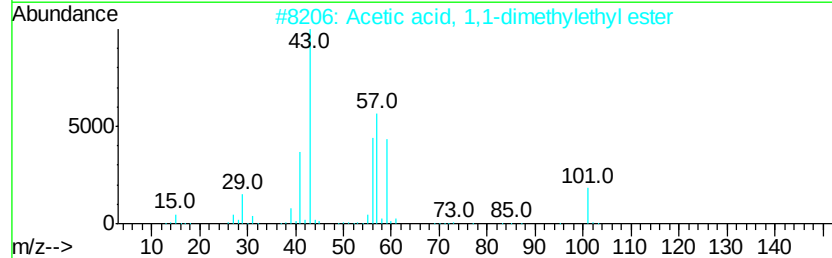
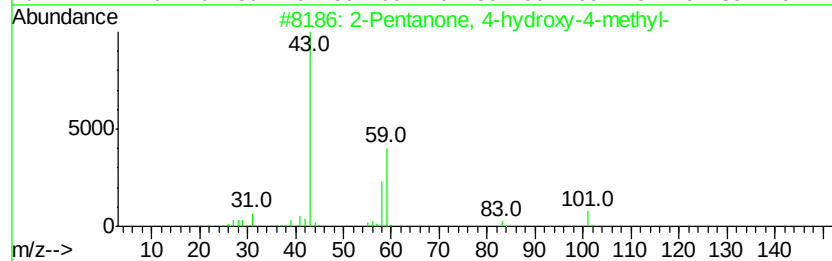
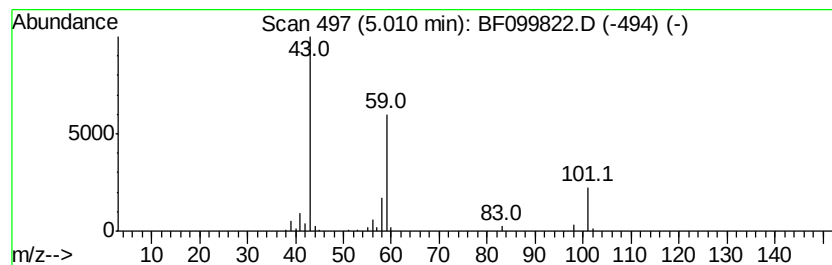
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.01	11.12 ng	212096	1,4-Dichlorobenzene-d4	6.79

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56
2			Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	39
3			Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	25
4			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9
5			Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9



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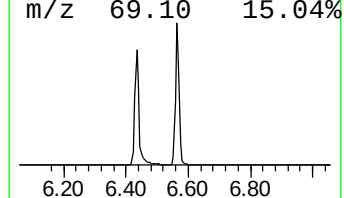
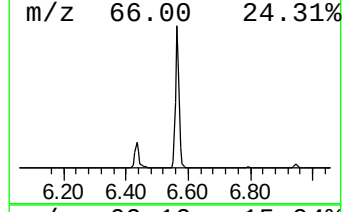
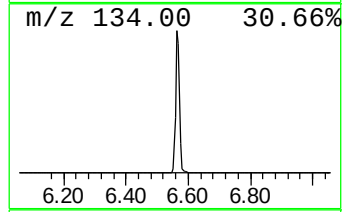
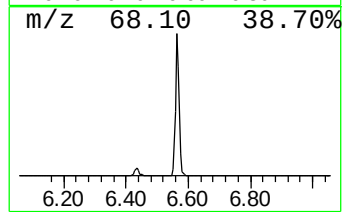
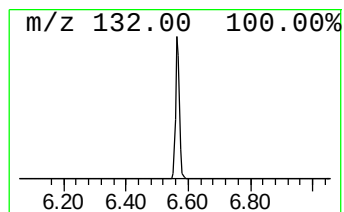
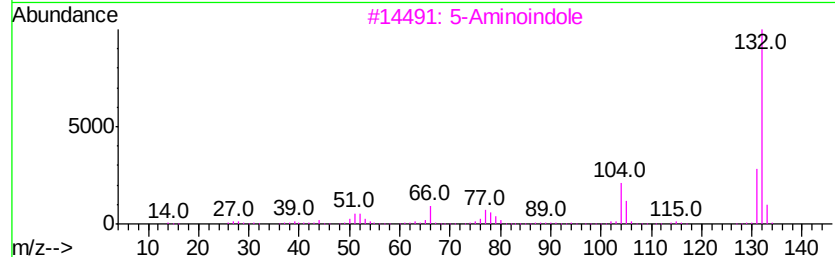
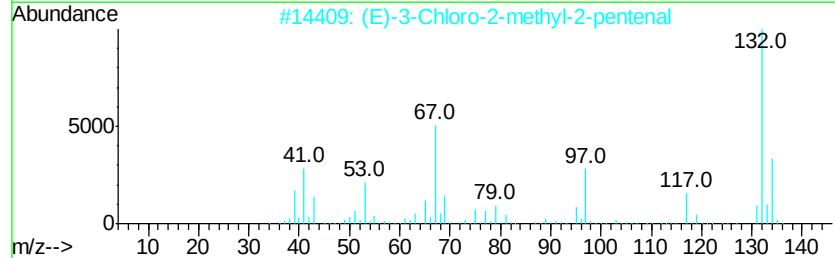
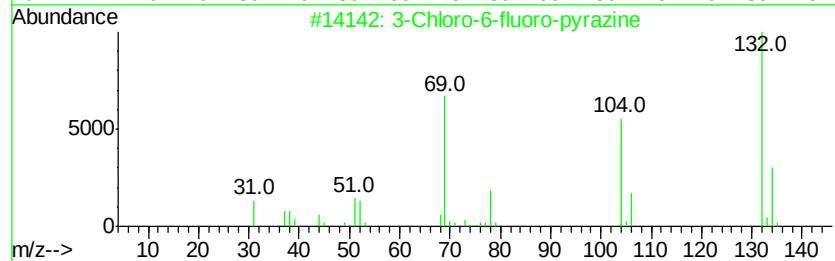
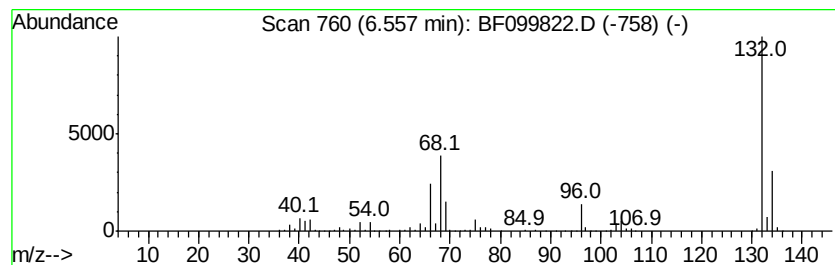
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Peak Number 2 unknown6.56 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.56	78.19 ng	1491130	1,4-Dichlorobenzene-d4	6.79

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	25
3		5-Aminoindole	132	C8H8N2	005192-03-0	12
4		Xenon	132	Xe	007440-63-3	9
5		1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.01	11.1	ng	212096	1	6.79	381390	20.0
unknown6.56	6.56	78.2	ng	1491130	1	6.79	381390	20.0