

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF100917\
 Data File : BF099261.D
 Acq On : 9 Oct 2017 14:42
 Operator : SJ/JU
 Sample : PB103065BL
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB103065BL

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-BF092317.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	5.104	509	513	525	rBV	403158	547506	16.29%	2.083%
2	5.498	575	580	583	rBV	2759629	2744860	81.65%	10.441%
3	6.498	745	750	753	rBV	2478112	2699554	80.30%	10.268%
4	6.634	769	773	777	rBV	2675289	2837538	84.41%	10.793%
5	6.863	808	812	815	rBV	744343	583517	17.36%	2.220%
6	7.022	834	839	842	rBV	2327539	2265100	67.38%	8.616%
7	7.428	902	908	911	rBV	1712907	1748395	52.01%	6.650%
8	8.145	1025	1030	1033	rBV	883321	800941	23.83%	3.047%
9	9.222	1207	1213	1216	rBV	3044706	3325838	98.93%	12.650%
10	9.898	1323	1328	1332	rBV2	1001697	890286	26.48%	3.386%
11	10.692	1457	1463	1466	rBV	1835408	1853221	55.13%	7.049%
12	11.386	1576	1581	1585	rBV	1085871	950310	28.27%	3.615%
13	12.974	1844	1851	1854	rBV	3199063	3361752	100.00%	12.787%
14	14.027	2026	2030	2037	rBV	1031557	894357	26.60%	3.402%
15	15.504	2275	2281	2292	rBV	603397	787214	23.42%	2.994%

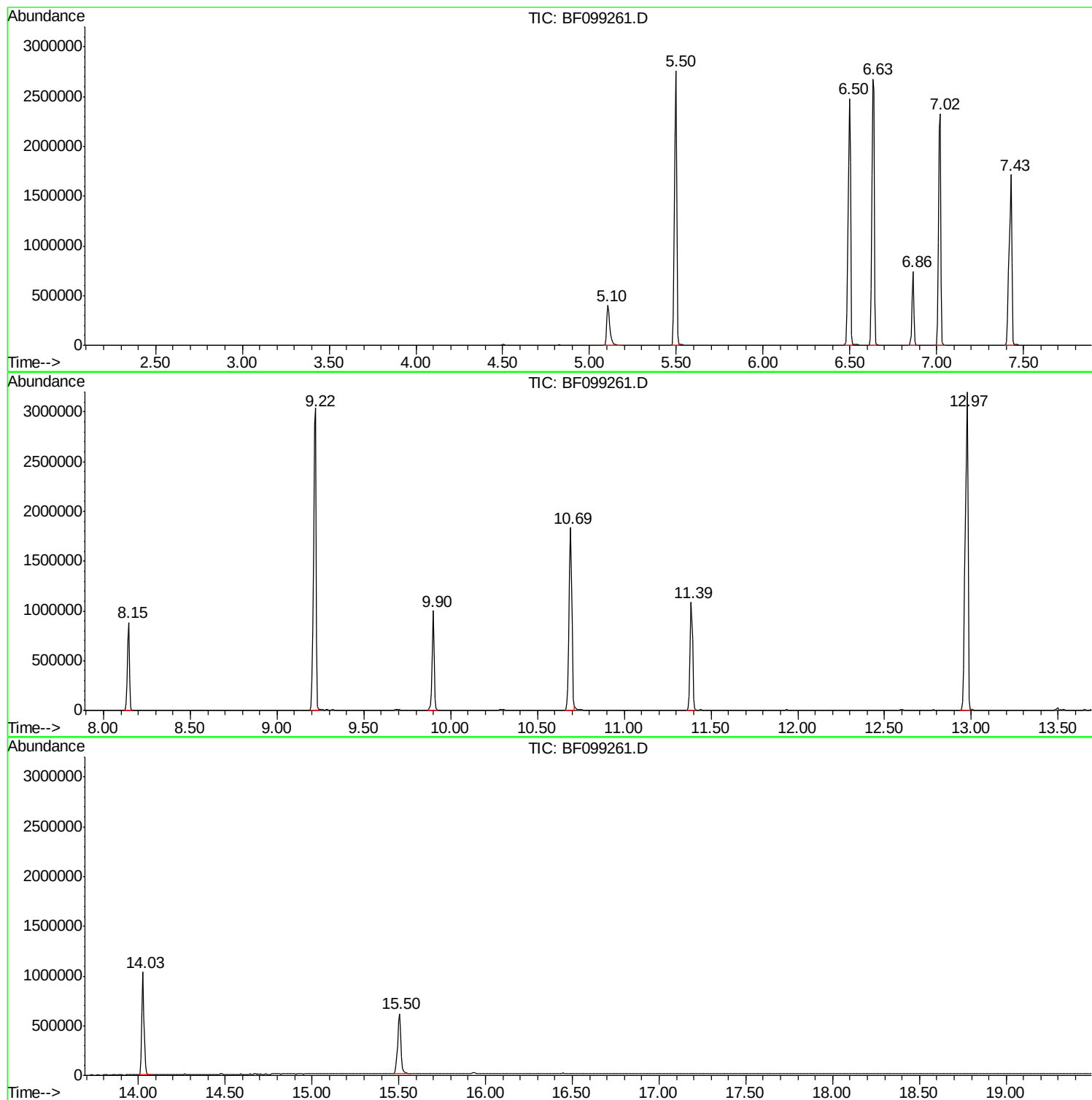
Sum of corrected areas: 26290389

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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF092317.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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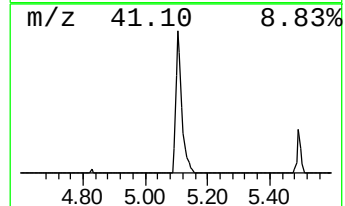
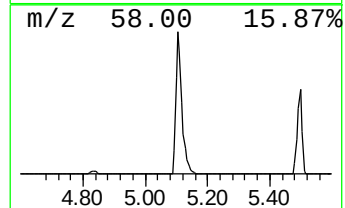
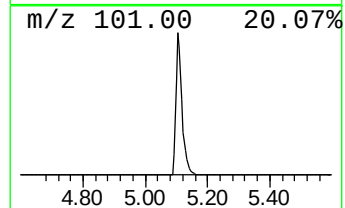
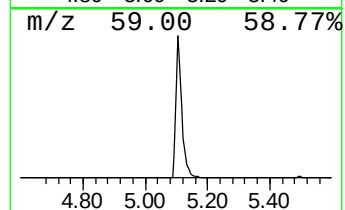
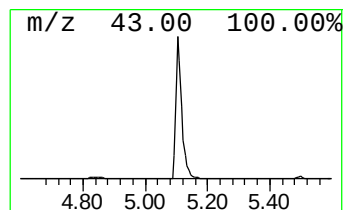
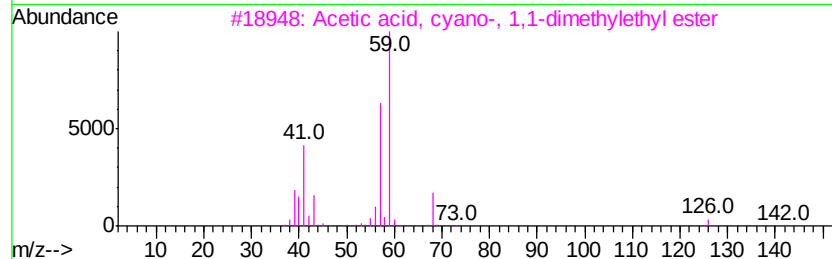
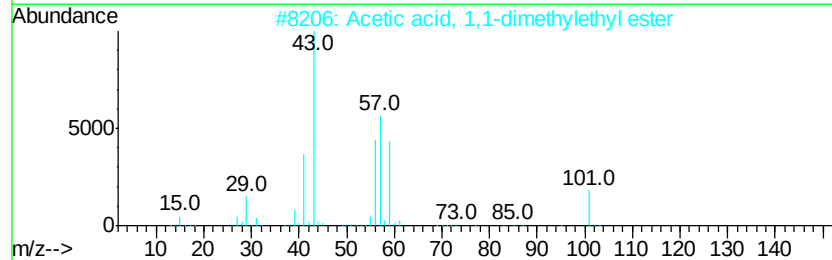
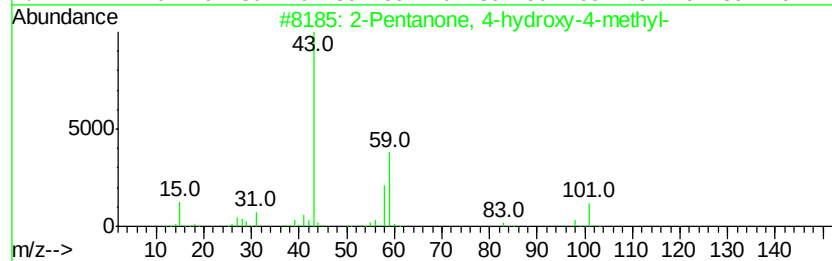
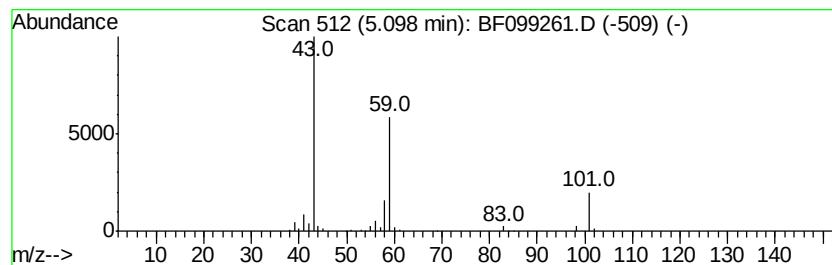
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.10	18.77 ng	547506	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	50
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	17
4			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9
5			Butane, 1-ethoxy-	102	C6H14O	000628-81-9	9



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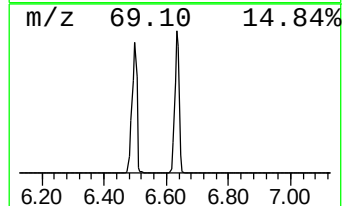
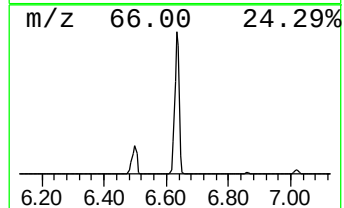
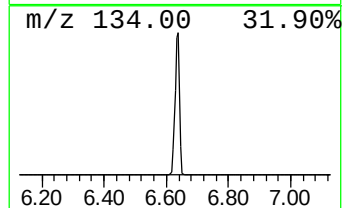
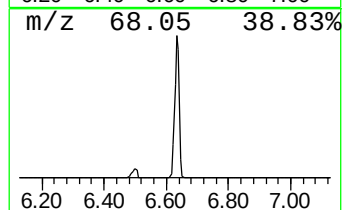
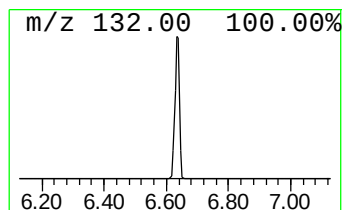
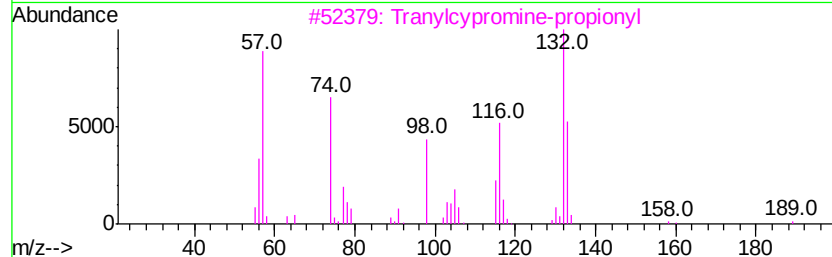
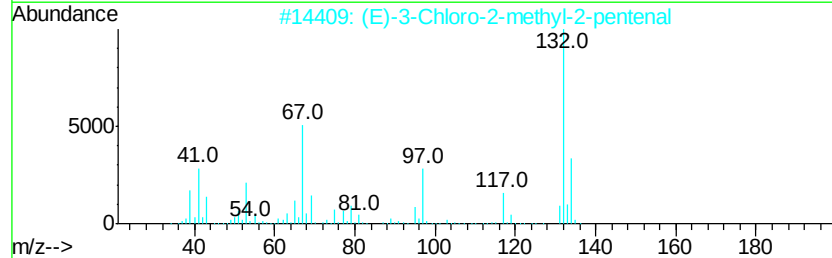
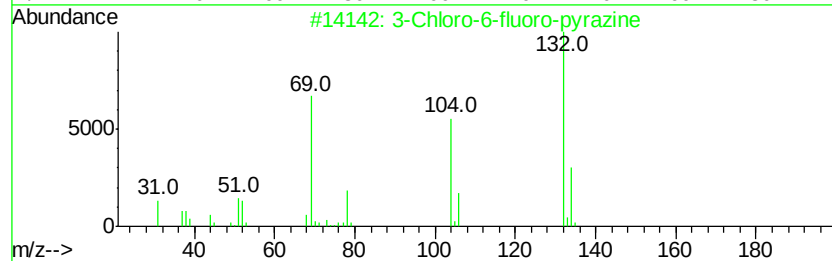
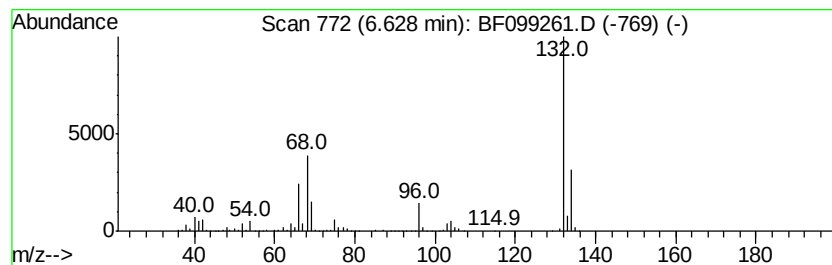
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Peak Number 2 unknown6.63 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.63	97.26 ng	2837540	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	35
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	17
3		Tranvlcypropine-propionyl	189	C12H15NO	1000123-86-3	12
4		5-Aminoindole	132	C8H8N2	005192-03-0	12
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...	5.10	18.8	ng	547506	1	6.86	583517	20.0
unknown6.63	6.63	97.3	ng	2837540	1	6.86	583517	20.0