

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF101117\
 Data File : BF099352.D
 Acq On : 11 Oct 2017 18:45
 Operator : SJ/JU
 Sample : PB103129BL
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB103129BL

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	5.092	507	511	527	rBV	296024	415441	13.90%	1.767%
2	5.481	573	577	591	rBV	2402888	2443743	81.79%	10.393%
3	6.487	742	748	751	rBV	2210392	2349835	78.64%	9.993%
4	6.622	766	771	774	rBV	2645490	2499543	83.66%	10.630%
5	6.851	806	810	813	rBV	743122	614612	20.57%	2.614%
6	7.004	832	836	840	rBV	2064308	2029434	67.92%	8.631%
7	7.416	900	906	909	rBV	1516717	1523146	50.98%	6.477%
8	8.133	1023	1028	1031	rBV	863613	868517	29.07%	3.694%
9	9.204	1204	1210	1214	rBV	2849393	2987917	100.00%	12.707%
10	9.886	1321	1326	1330	rBV	1079588	964571	32.28%	4.102%
11	10.674	1455	1460	1464	rBV	1531019	1602306	53.63%	6.814%
12	11.374	1574	1579	1582	rBV	1152499	987151	33.04%	4.198%
13	12.957	1843	1848	1851	rBV	2859944	2785748	93.23%	11.847%
14	13.480	1932	1937	1941	rBV	36569	39911	1.34%	0.170%
15	14.010	2023	2027	2031	rBV	933656	821785	27.50%	3.495%
16	15.480	2272	2277	2284	rBV	461178	580783	19.44%	2.470%

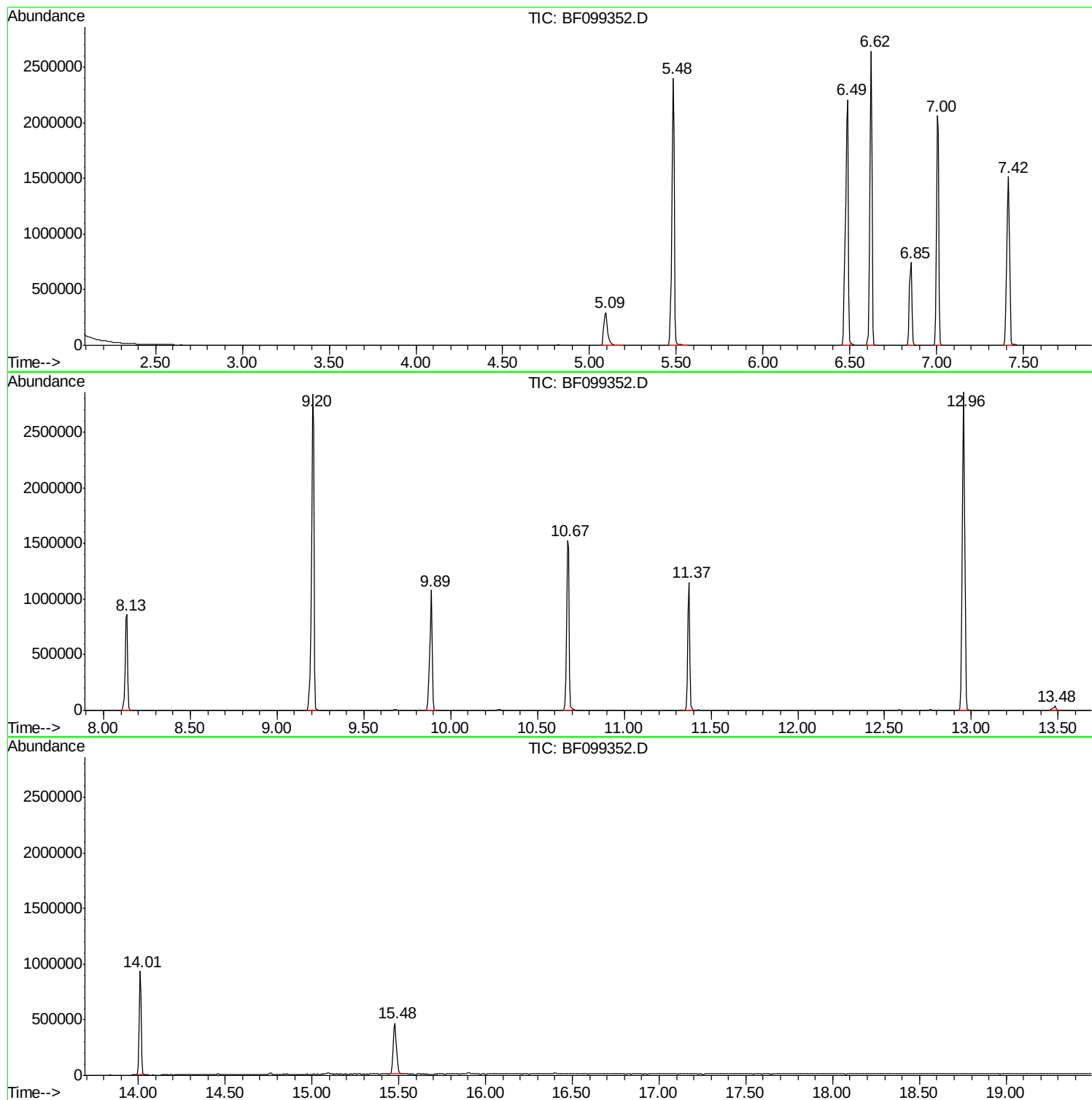
Sum of corrected areas: 23514443

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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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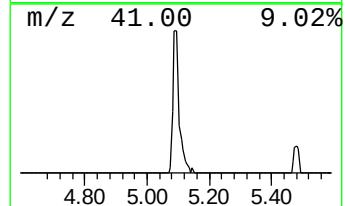
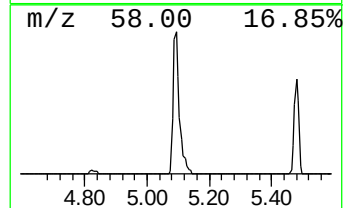
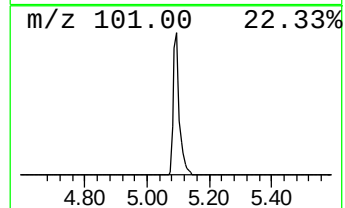
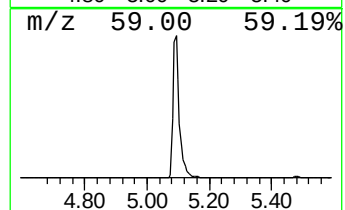
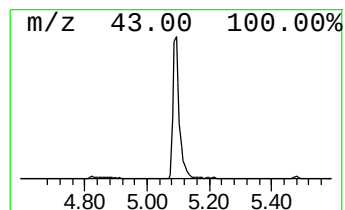
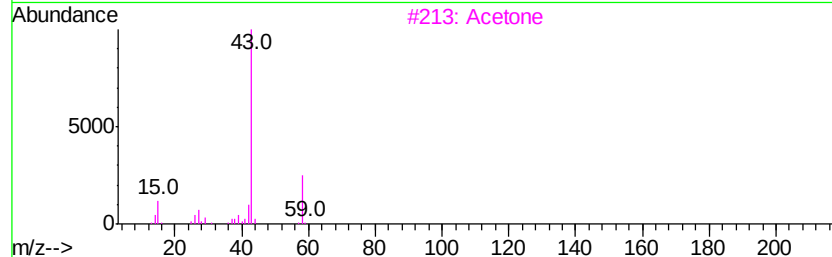
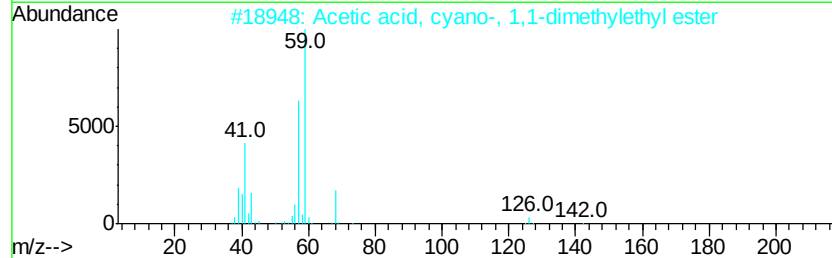
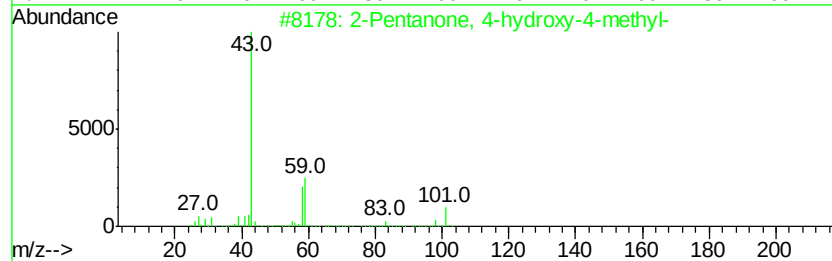
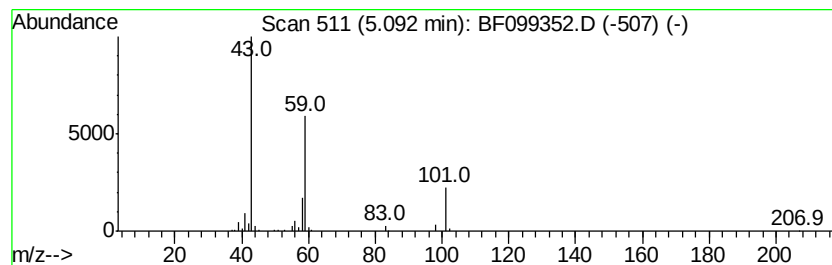
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.09	13.52 ng	415441	1,4-Dichlorobenzene-d4	6.85

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	64
2		Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	23
3		Acetone	58	C3H6O	000067-64-1	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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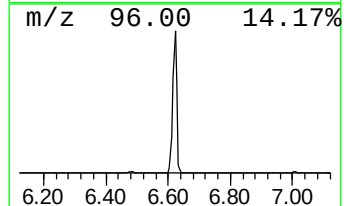
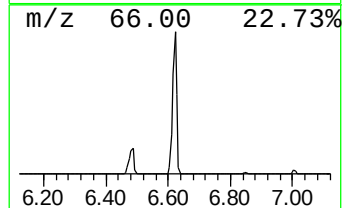
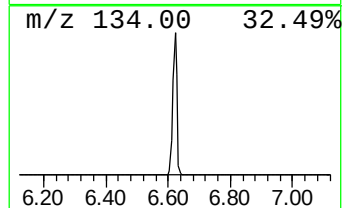
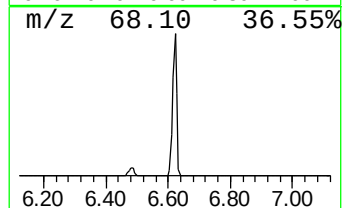
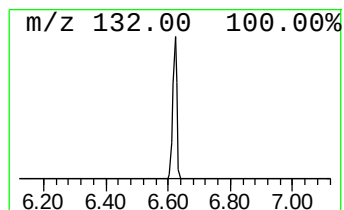
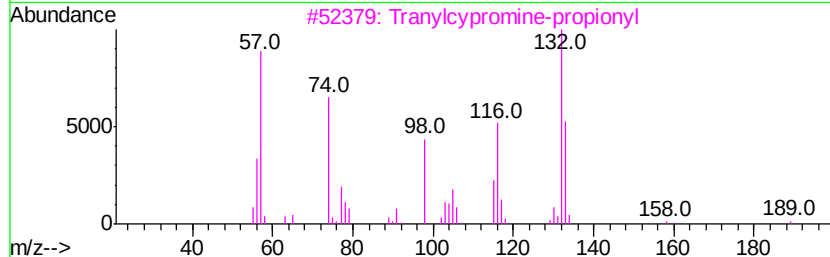
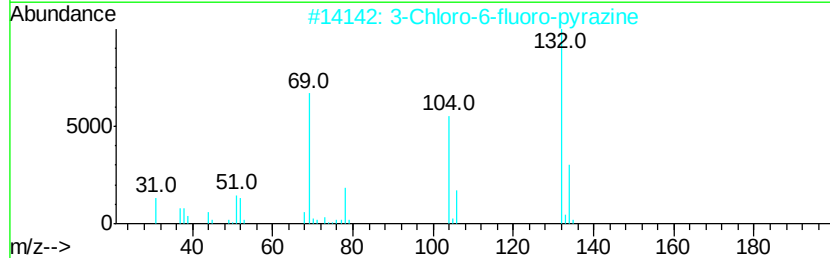
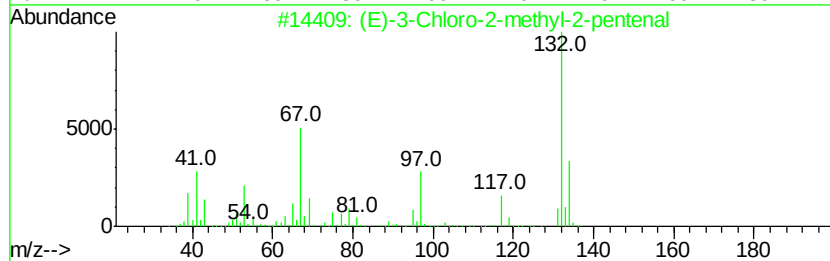
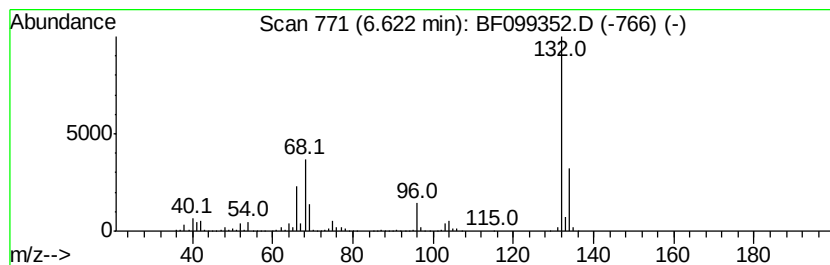
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Peak Number 2 unknown6.62 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.62	81.34 ng	2499540	1,4-Dichlorobenzene-d4	6.85

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	35
2		3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	25
3		Tranvlcypromine-propionyl	189	C12H15NO	1000123-86-3	12
4		5-Aminoindole	132	C8H8N2	005192-03-0	12
5		4-Chloro-2-fluoro-pyrimidine	132	C4H2ClFN2	051422-00-5	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.09	13.5	ng	415441	1	6.85	614612	20.0
unknown6.62	6.62	81.3	ng	2499540	1	6.85	614612	20.0