

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF111717\
 Data File : BF100674.D
 Acq On : 17 Nov 2017 14:24
 Operator : SJ/JU
 Sample : I6454-04
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 N-NARANJITO-S001-0001-02

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-BF110817.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	3.481	235	237	247	rBV3	14393	39582	1.19%	0.149%
2	5.116	511	515	531	rBB	358367	553034	16.69%	2.076%
3	5.510	576	582	585	rBV	2664582	2942114	88.80%	11.046%
4	6.516	746	753	755	rBV	2475585	3081095	92.99%	11.568%
5	6.657	771	777	780	rBV	2699370	3036343	91.64%	11.399%
6	6.880	811	815	819	rVB	741525	596861	18.01%	2.241%
7	7.039	837	842	845	rVB	2471746	2316980	69.93%	8.699%
8	7.445	905	911	914	rBV	1795631	1993757	60.17%	7.485%
9	7.516	919	923	931	rVB	35176	35246	1.06%	0.132%
10	8.163	1028	1033	1036	rBV	899700	780826	23.57%	2.931%
11	9.239	1210	1216	1219	rBV	3312308	3313271	100.00%	12.439%
12	9.627	1277	1282	1285	rBV	552682	438869	13.25%	1.648%
13	9.915	1326	1331	1335	rBV	979000	862978	26.05%	3.240%
14	10.709	1458	1466	1469	rBV	1798945	1871306	56.48%	7.026%
15	10.809	1480	1483	1493	rVB	42151	46474	1.40%	0.174%
16	11.404	1580	1584	1587	rBV	1014560	862762	26.04%	3.239%
17	12.986	1848	1853	1857	rBV	2534134	2608950	78.74%	9.795%
18	13.868	2000	2003	2012	rBV2	45150	52815	1.59%	0.198%
19	14.039	2028	2032	2035	rBV	683246	585547	17.67%	2.198%
20	15.515	2277	2283	2292	rBV	454746	616959	18.62%	2.316%

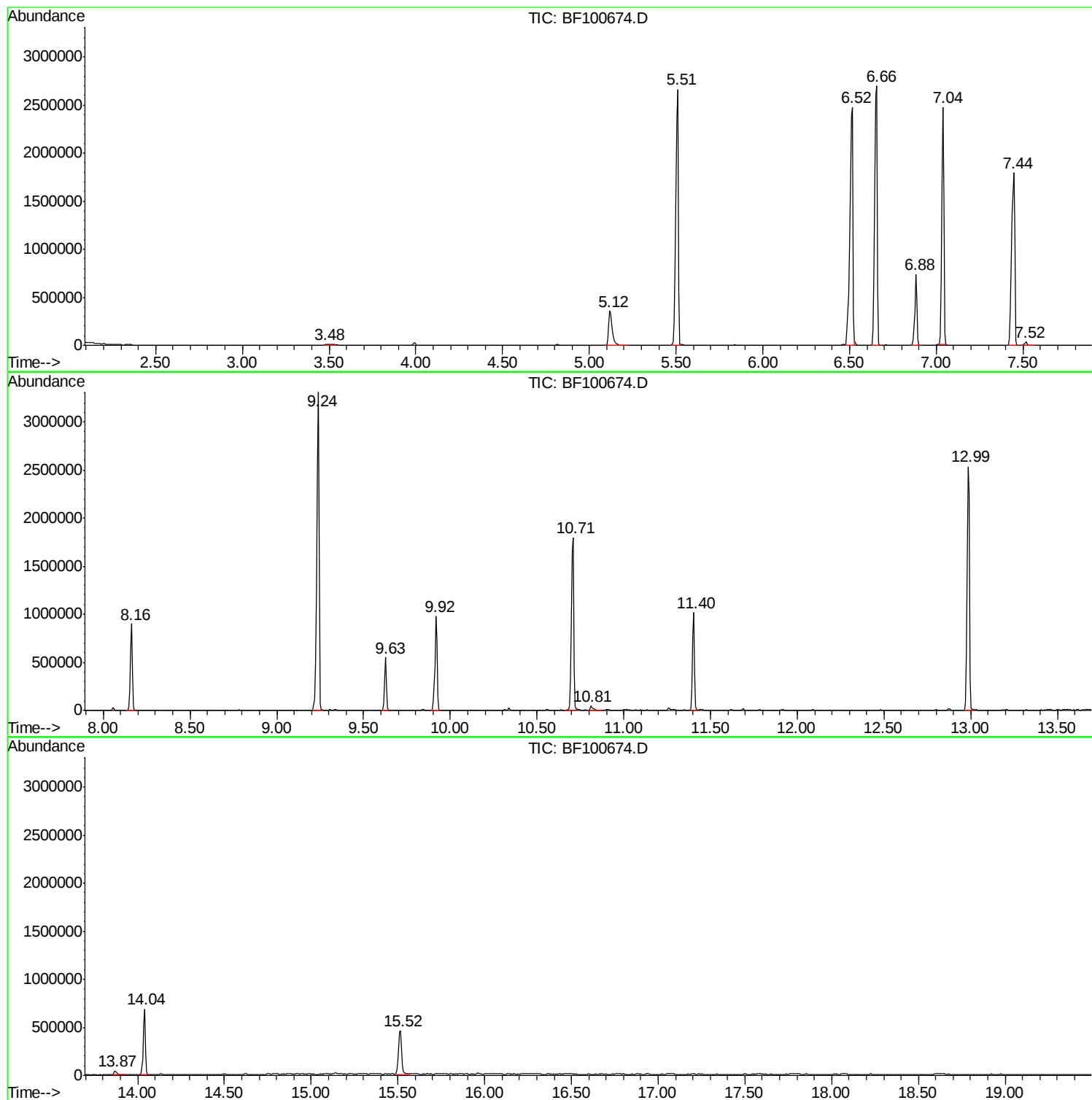
Sum of corrected areas: 26635769

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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF110817.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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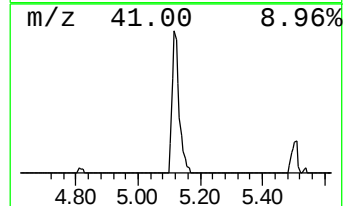
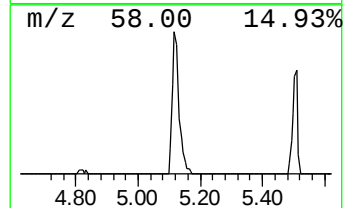
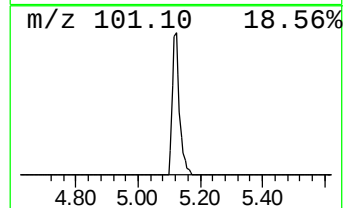
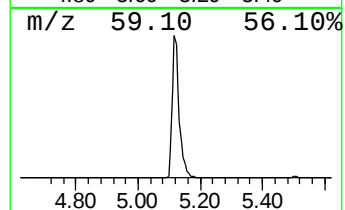
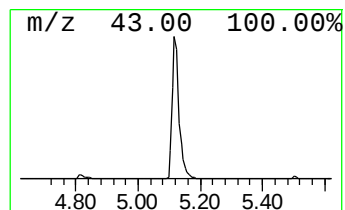
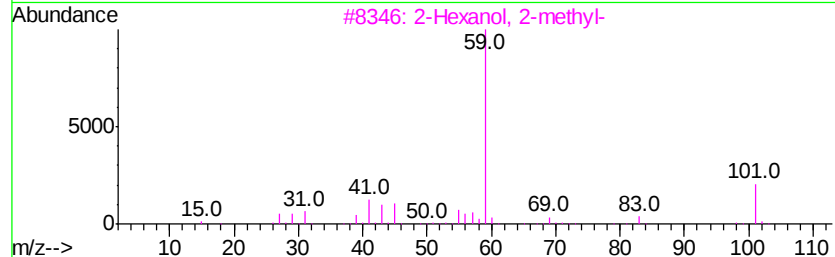
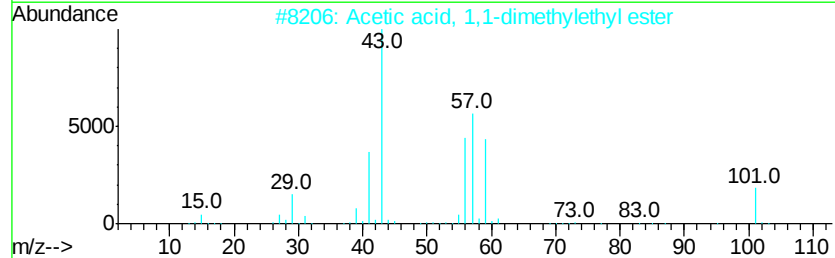
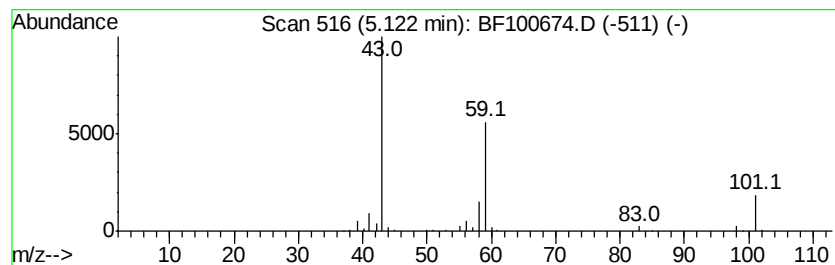
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.12	18.53 ng	553034	1,4-Dichlorobenzene-d4	6.88

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	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		2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	28
4		Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	23
5		Formamide, N,N-diethyl-	101	C5H11NO	000617-84-5	9



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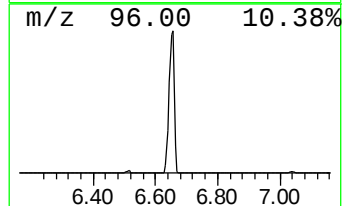
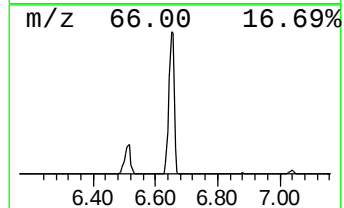
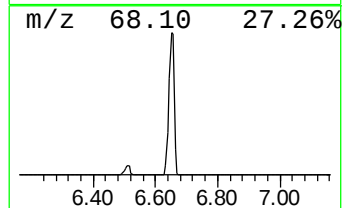
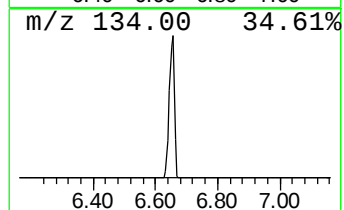
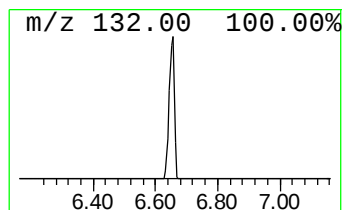
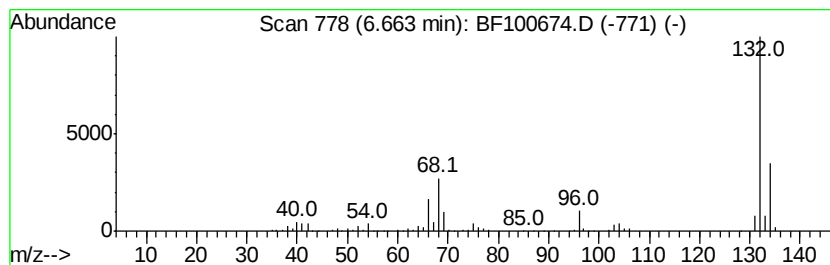
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Peak Number 2 unknown6.66 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.66	101.74 ng	3036340	1,4-Dichlorobenzene-d4	6.88

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	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	5	Aminoindole	132	C8H8N2	005192-03-0	12
4	Imidazole, 4,5	dicvano-1-methyl-	132	C6H4N4	019485-35-9	9
5	1,3,5	Trifluorobenzene	132	C6H3F3	000372-38-3	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.12	18.5	ng	553034	1	6.88	596861	20.0
unknown6.66	6.66	101.7	ng	3036340	1	6.88	596861	20.0