

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF111516\
 Data File : BF091179.D
 Acq On : 15 Nov 2016 15:26
 Operator : UM/SJ
 Sample : H5636-06
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 C19043041

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-BF110316.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	1.744	3	5	58	rVB	3107396	9477736	100.00%	18.640%
2	4.762	267	269	280	rBV	365326	515199	5.44%	1.013%
3	5.219	307	309	326	rBV	1978502	2301335	24.28%	4.526%
4	6.282	400	402	410	rBV	1250963	1498069	15.81%	2.946%
5	6.407	410	413	415	rBV	3186322	4295842	45.33%	8.449%
6	6.636	431	433	442	rBV	884397	999950	10.55%	1.967%
7	6.785	444	446	449	rBV	3387682	4193504	44.25%	8.247%
8	7.207	480	483	485	rBV	3236906	3389532	35.76%	6.666%
9	7.916	543	545	548	rBV	1082977	1340971	14.15%	2.637%
10	8.339	579	582	590	rBV	61301	132834	1.40%	0.261%
11	9.002	637	640	642	rBV	5814976	5959526	62.88%	11.720%
12	9.402	673	675	680	rBV	169820	162980	1.72%	0.321%
13	9.665	696	698	701	rBV	1573766	1658133	17.50%	3.261%
14	10.099	731	736	739	rBV4	77946	166692	1.76%	0.328%
15	10.465	765	768	770	rBV	3084599	3781290	39.90%	7.437%
16	11.151	826	828	830	rBV	1627789	1629328	17.19%	3.204%
17	12.739	964	967	969	rBV	5775212	6458612	68.15%	12.702%
18	13.780	1055	1058	1060	rBV	1493444	1603118	16.91%	3.153%
19	15.140	1175	1177	1180	rBV	848512	1282661	13.53%	2.523%

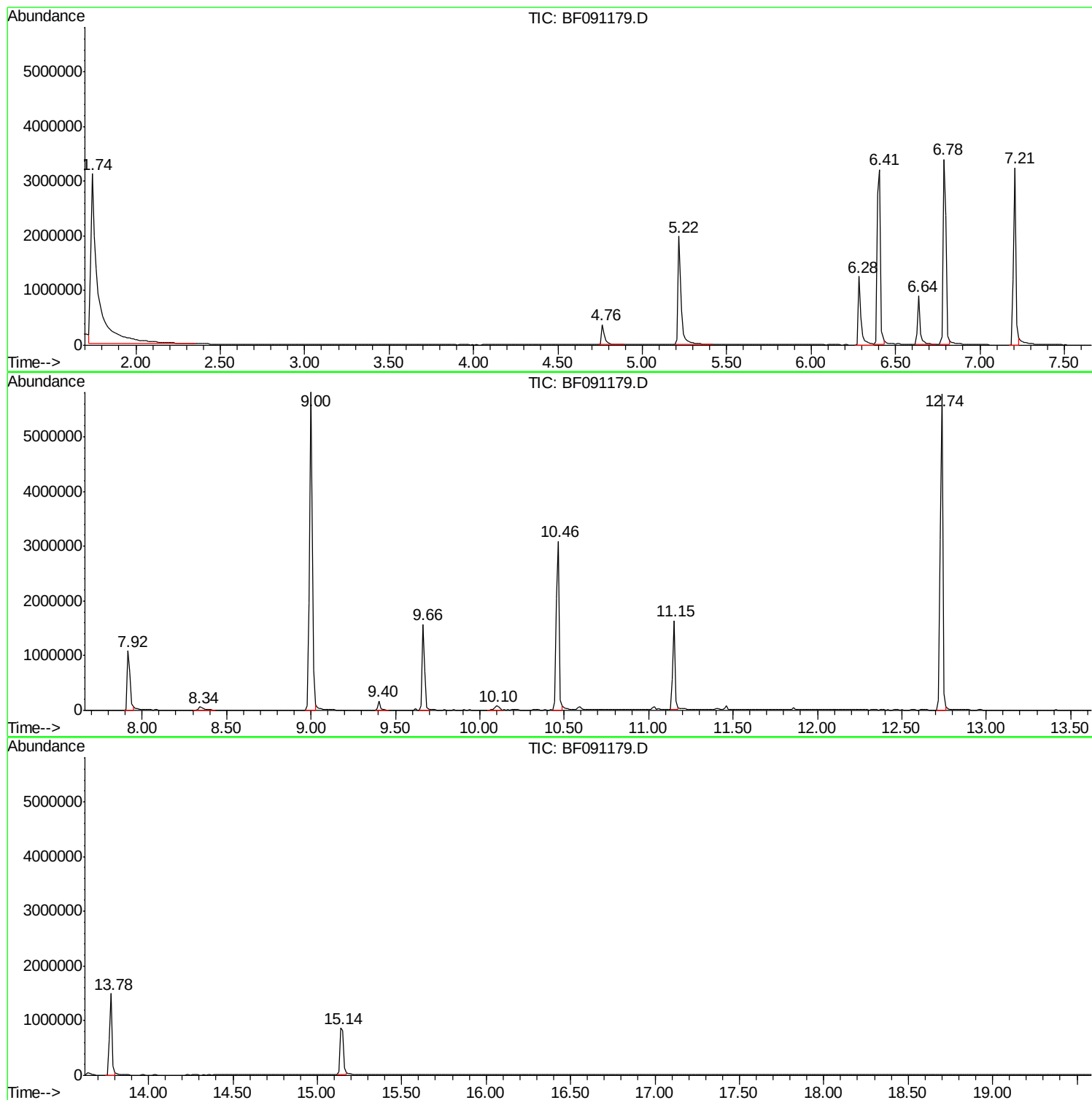
Sum of corrected areas: 50847312

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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF110316.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P



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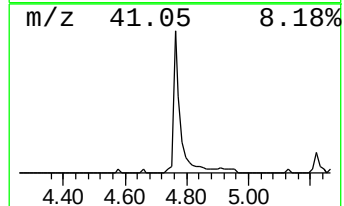
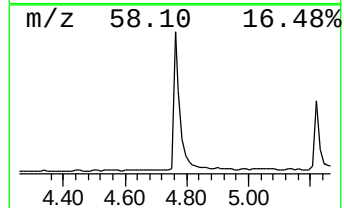
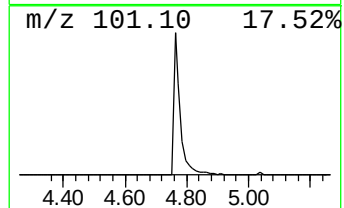
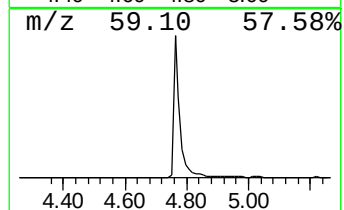
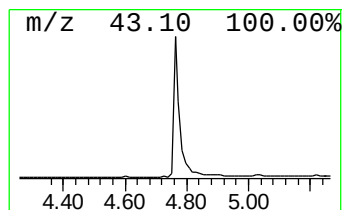
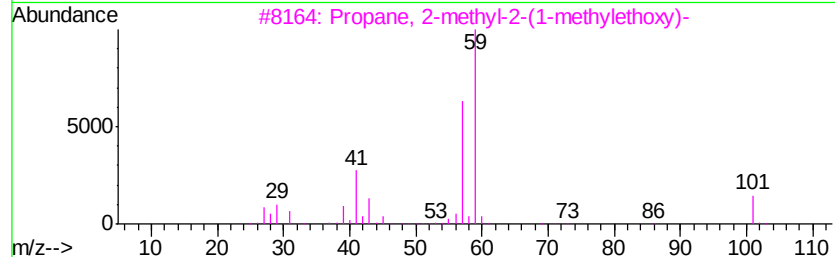
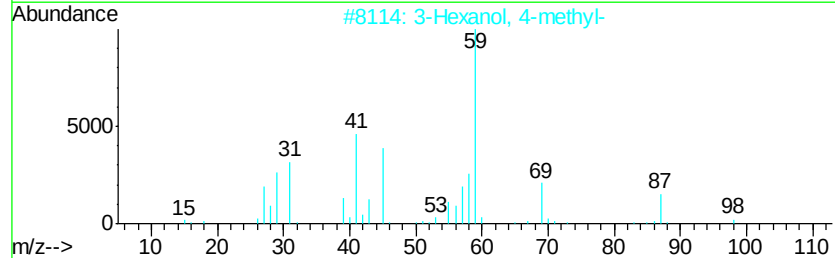
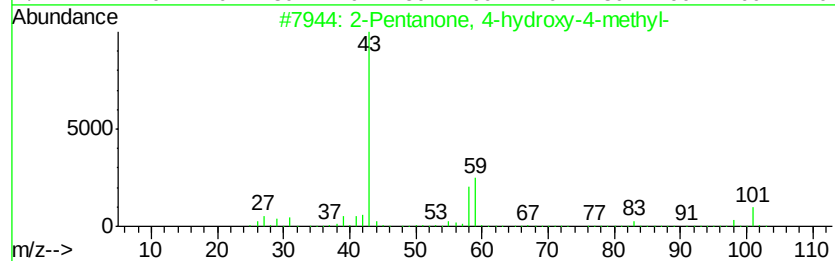
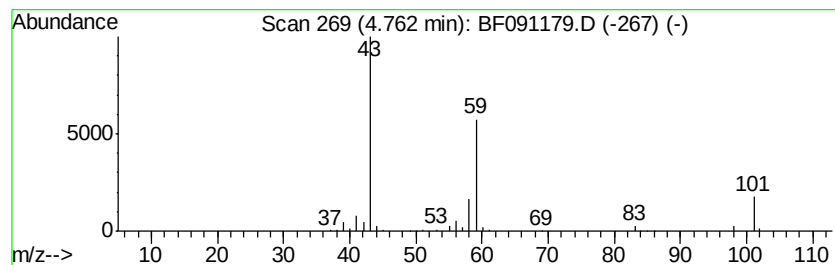
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Peak Number 2 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.76	10.30 ng	515199	1,4-Dichlorobenzene-d4	6.64

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	59
2			3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	33
3			Propane, 2-methyl-2-(1-methylethoxy)-	116	C7H16O	017348-59-3	33
4			2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	28
5			Acetic acid, cyano-, 1,1-dimethyl-	141	C7H11NO2	001116-98-9	25



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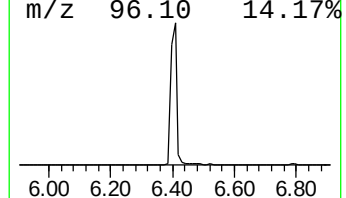
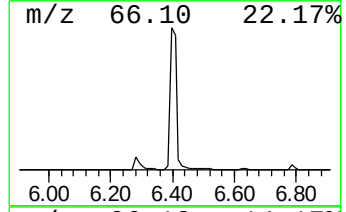
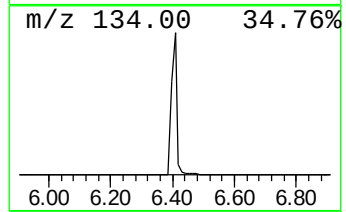
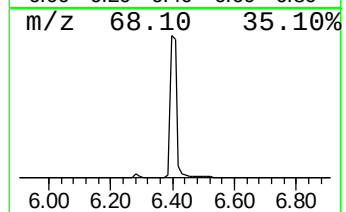
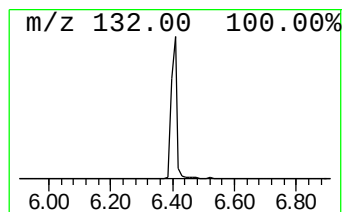
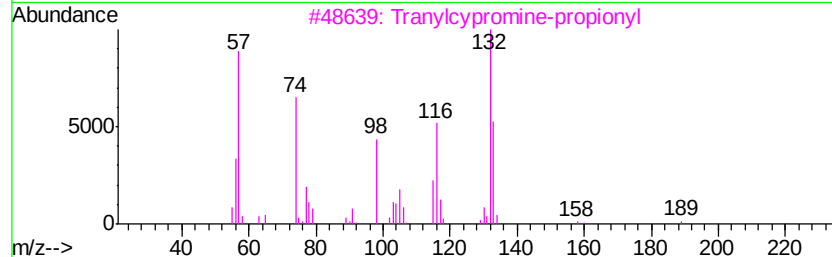
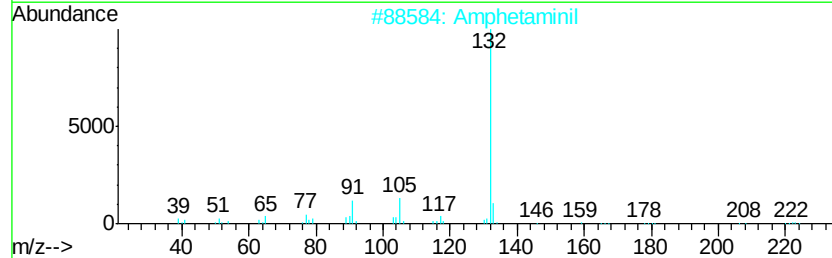
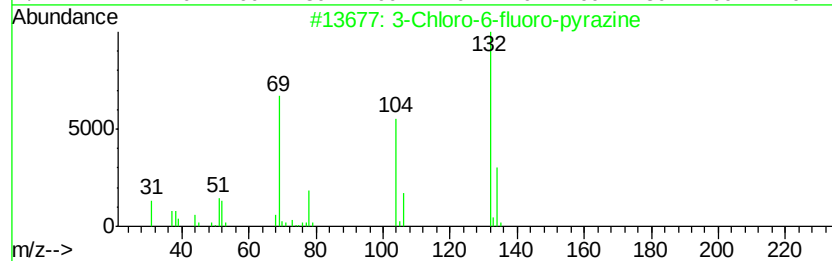
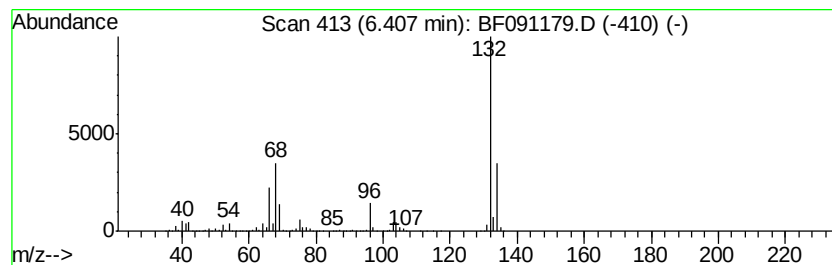
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Peak Number 3 unknown6.41 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.41	85.92 ng	4295840	1,4-Dichlorobenzene-d4	6.64

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	25
2		Amphetaminil	250	C17H18N2	017590-01-1	12
3		Tranvlcyproline-propionyl	189	C12H15NO	1000123-86-3	12
4		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	12
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...	4.76	10.3	ng	515199	1	6.64	999950	20.0
unknown6.41	6.41	85.9	ng	4295840	1	6.64	999950	20.0