

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF062216\
 Data File : BF088257.D
 Acq On : 22 Jun 2016 18:50
 Operator : UM/SJ
 Sample : H3660-09
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 GW-1

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-BF060716.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	1.724	11	12	56	rVB	1564064	3438277	100.00%	29.648%
2	4.719	272	274	291	rBV	152935	186977	5.44%	1.612%
3	5.199	314	316	331	rBV	98652	158270	4.60%	1.365%
4	6.262	407	409	417	rBV	99711	122649	3.57%	1.058%
5	6.376	417	419	424	rVB	227331	315424	9.17%	2.720%
6	6.604	437	439	446	rBB	378590	378172	11.00%	3.261%
7	6.753	450	452	455	rBV	753923	685454	19.94%	5.911%
8	7.176	487	489	491	rBV	449961	475707	13.84%	4.102%
9	7.885	550	551	554	rBV	555565	591449	17.20%	5.100%
10	8.970	644	646	648	rBV	1018933	1121985	32.63%	9.675%
11	9.370	679	681	683	rBV	133551	105036	3.05%	0.906%
12	9.633	702	704	707	rBV	734636	684361	19.90%	5.901%
13	10.079	741	743	745	rVB	56529	44304	1.29%	0.382%
14	10.422	771	773	776	rBV	252197	233971	6.80%	2.017%
15	10.548	783	784	789	rVB	52946	57190	1.66%	0.493%
16	11.119	832	834	836	rBV	497845	671407	19.53%	5.789%
17	12.525	955	957	959	rBV	45902	42298	1.23%	0.365%
18	12.697	970	972	974	rBV	1088610	1098278	31.94%	9.470%
19	13.622	1049	1053	1061	rBV	20441	62567	1.82%	0.540%
20	13.748	1061	1064	1070	rVB	430050	657225	19.11%	5.667%
21	15.108	1181	1183	1191	rBV	346944	466142	13.56%	4.019%

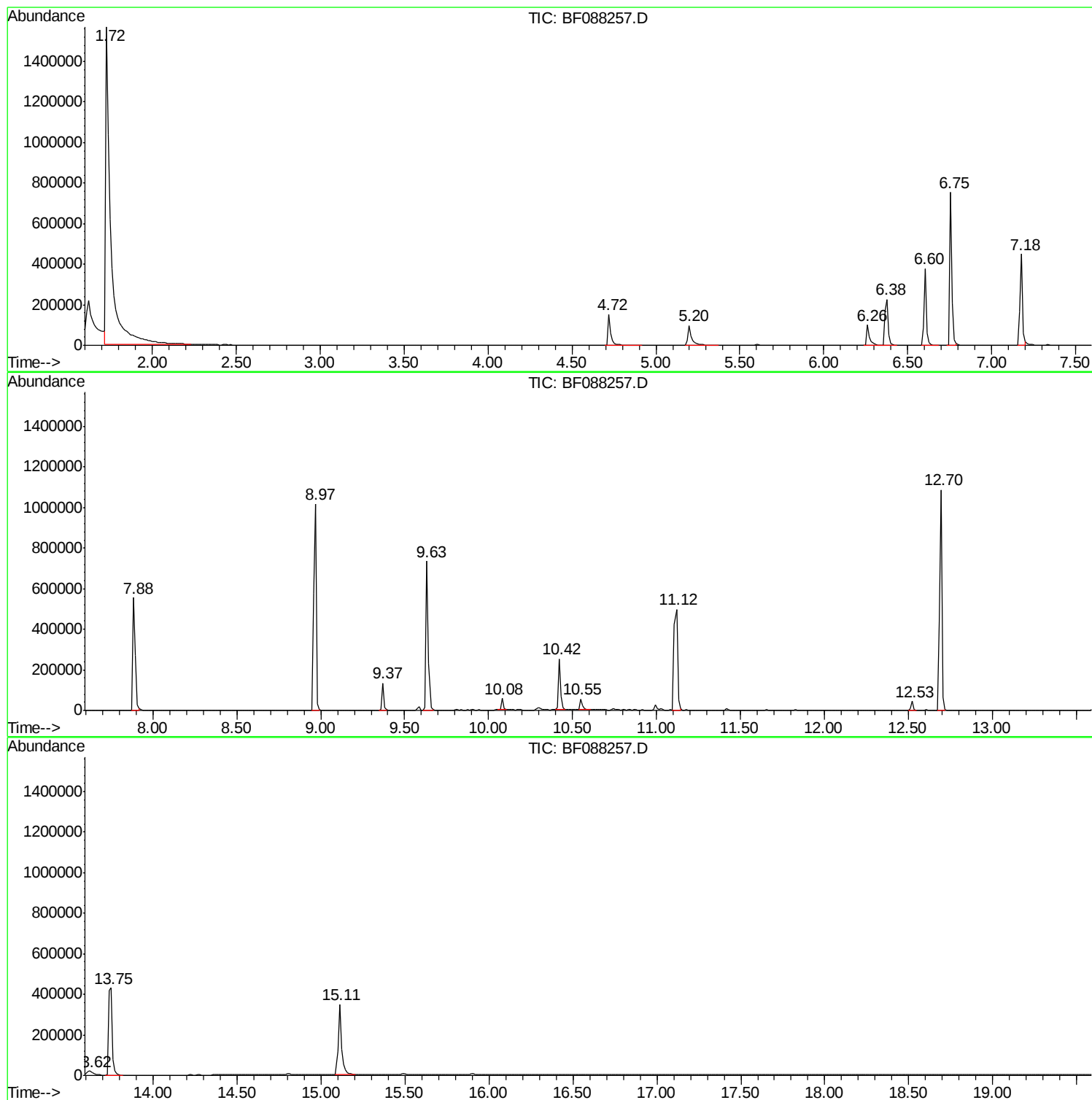
Sum of corrected areas: 11597143

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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF060716.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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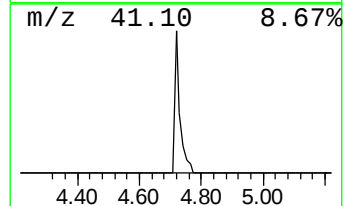
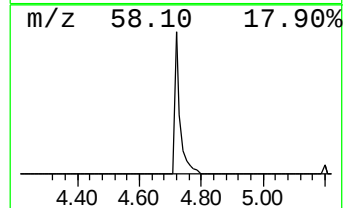
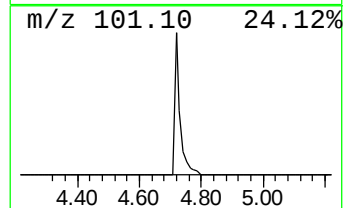
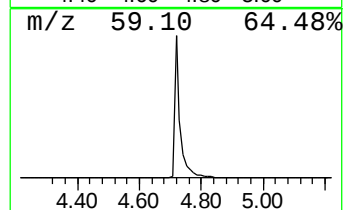
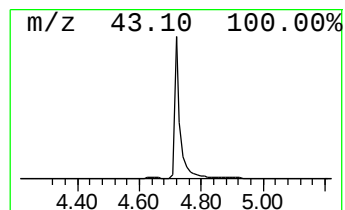
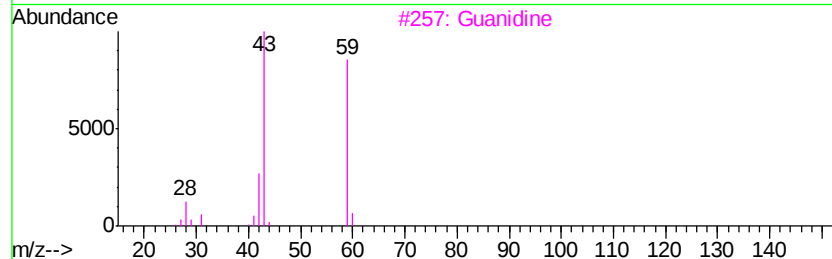
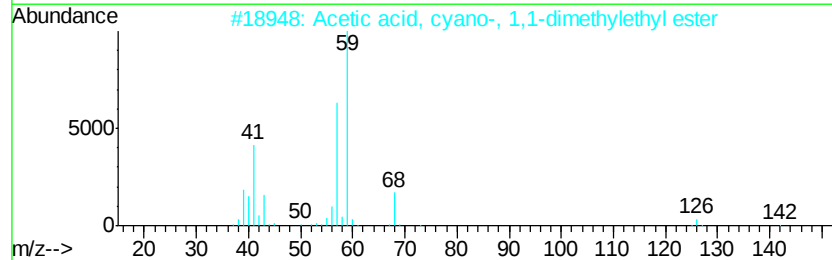
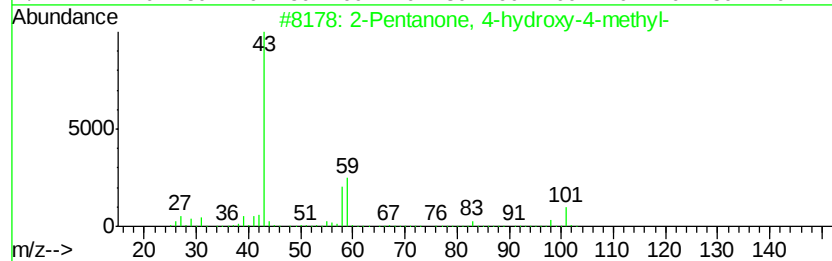
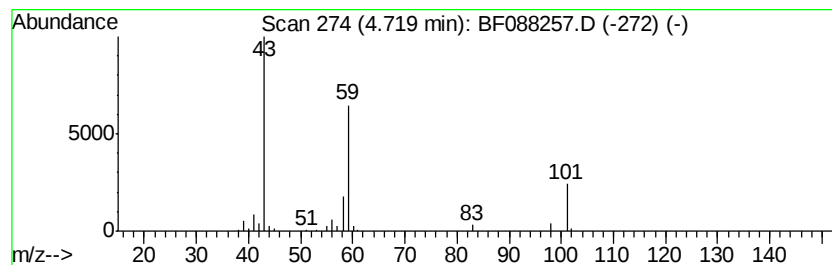
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Peak Number 2 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.72	9.89 ng	186977	1,4-Dichlorobenzene-d4	6.60

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	59
2		Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	17
3		Guanidine	59	CH5N3	000113-00-8	9
4		Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9
5		Acetone	58	C3H6O	000067-64-1	9



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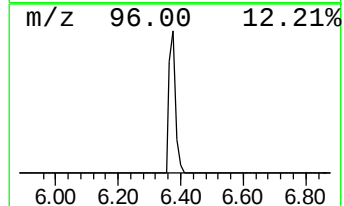
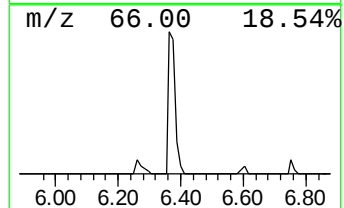
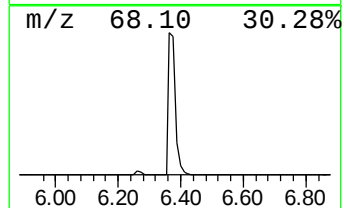
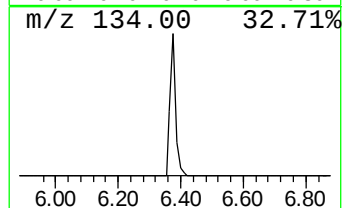
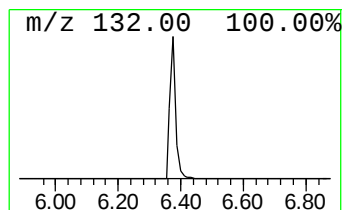
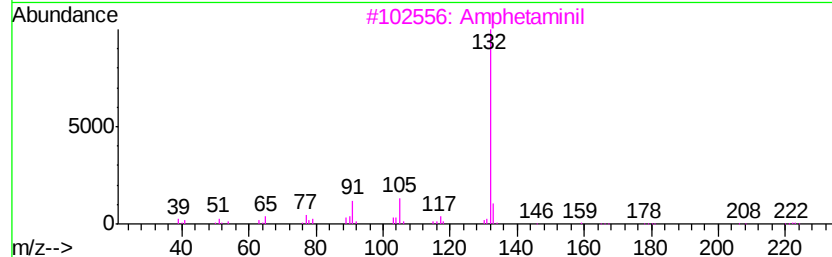
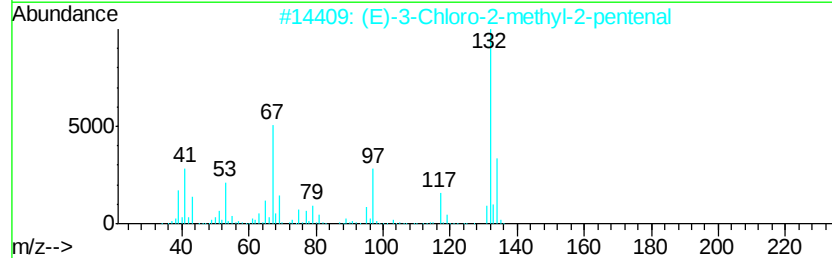
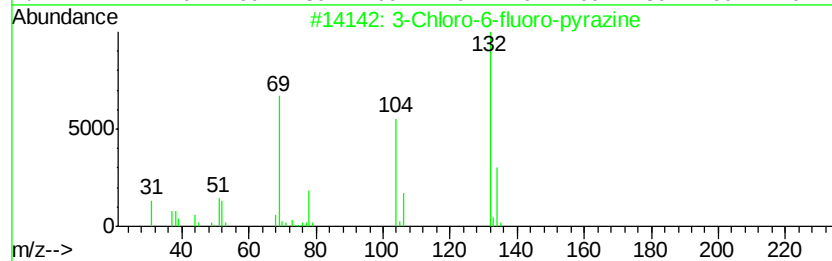
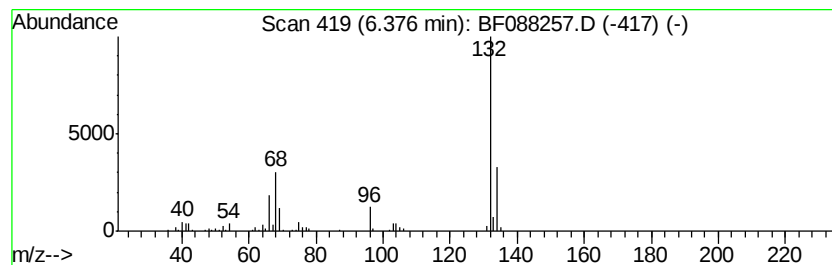
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Peak Number 3 unknown6.38 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.38	16.68 ng	315424	1,4-Dichlorobenzene-d4	6.60

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	40
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	33
3		Amphetaminil	250	C17H18N2	017590-01-1	25
4		4-Chloro-6-fluoro-pyrimidine	132	C4H2ClFN2	051422-01-6	9
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...	4.72	9.9	ng	186977	1	6.60	378172	20.0
unknown6.38	6.38	16.7	ng	315424	1	6.60	378172	20.0