

Data Path : Z:\HPCHEM1\BNA F\DATA\BF061516\
 Data File : BF088042.D
 Acq On : 15 Jun 2016 19:10
 Operator : UM/SJ
 Sample : H3609-01
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 181-01-SW-06142016

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	1.701	7	10	18	rVB	204334	470521	15.83%	2.333%
2	1.815	18	20	31	rVB	1719516	2973168	100.00%	14.742%
3	4.901	288	290	299	rBV	158081	195092	6.56%	0.967%
4	5.336	325	328	330	rBV	1513255	1615114	54.32%	8.008%
5	6.376	417	419	422	rBV	1193960	1543004	51.90%	7.651%
6	6.513	428	431	433	rVB	1217436	1666033	56.04%	8.261%
7	6.741	448	451	453	rBV	709299	595193	20.02%	2.951%
8	6.901	462	465	467	rBV	1151457	1398129	47.02%	6.933%
9	7.313	498	501	503	rBV	765782	1071490	36.04%	5.313%
10	8.033	561	564	566	rBV	565209	769605	25.89%	3.816%
11	9.107	655	658	660	rBV	2105200	1952128	65.66%	9.680%
12	9.507	690	693	695	rBV	250573	273231	9.19%	1.355%
13	9.782	714	717	719	rVB	892891	861837	28.99%	4.273%
14	10.216	754	755	757	rVB	32253	30363	1.02%	0.151%
15	10.570	783	786	788	rBV2	729166	1002533	33.72%	4.971%
16	10.696	795	797	801	rBV	29542	46322	1.56%	0.230%
17	11.142	834	836	837	rBV	27805	30064	1.01%	0.149%
18	11.256	844	846	849	rBV	886912	807138	27.15%	4.002%
19	12.845	982	985	987	rBV	1796991	1636341	55.04%	8.114%
20	13.885	1074	1076	1079	rBV2	525804	653704	21.99%	3.241%
21	15.291	1196	1199	1205	rVB	448618	576556	19.39%	2.859%

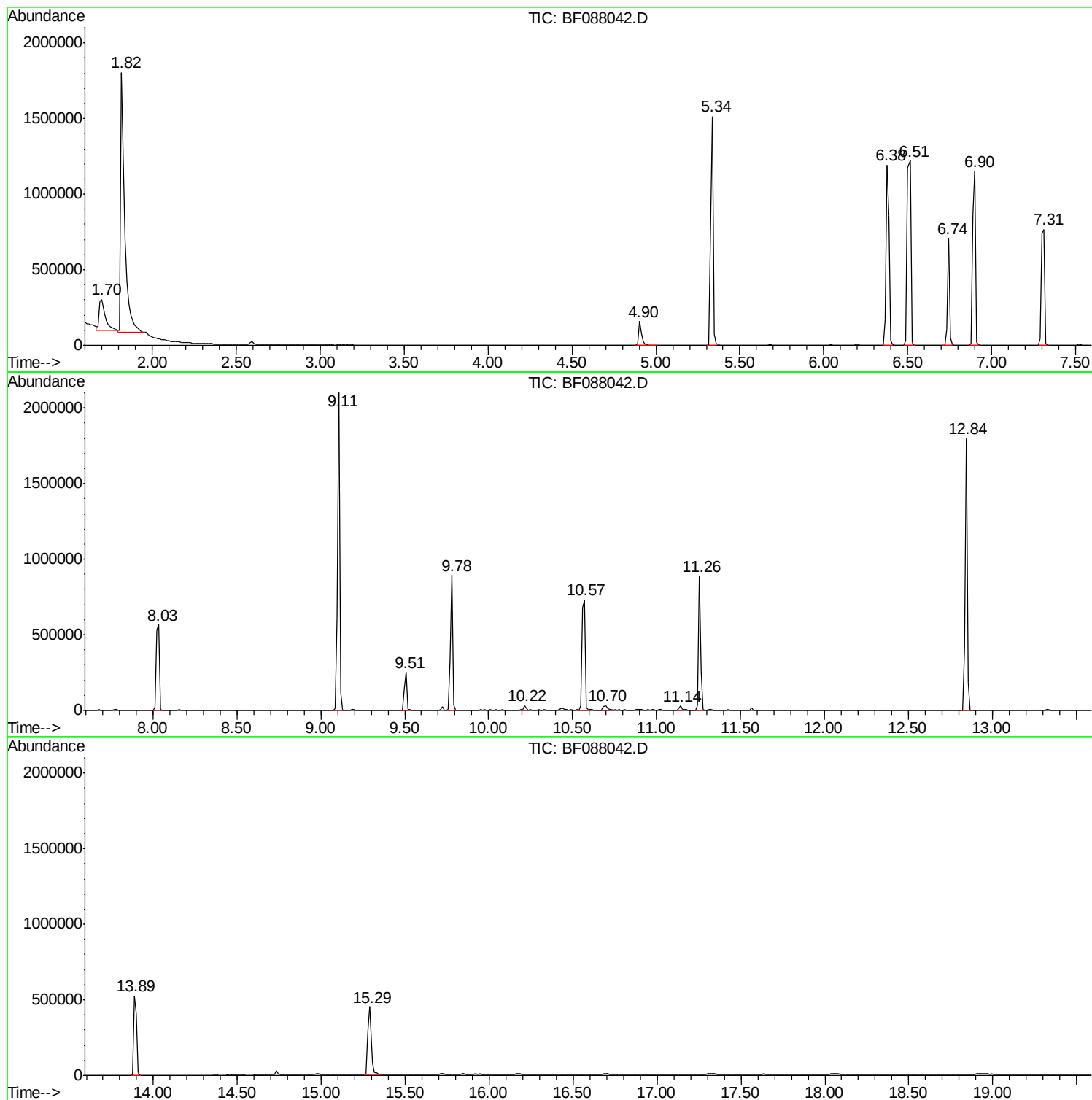
Sum of corrected areas: 20167566

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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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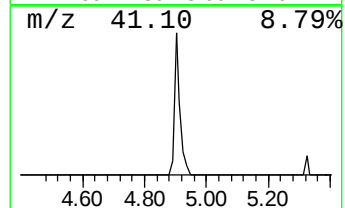
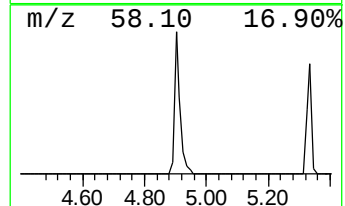
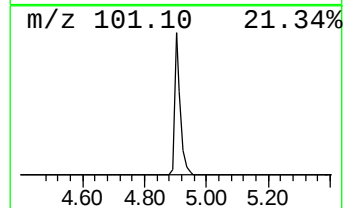
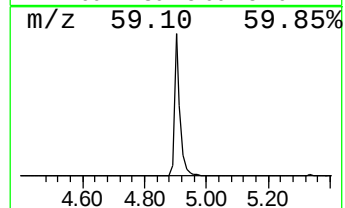
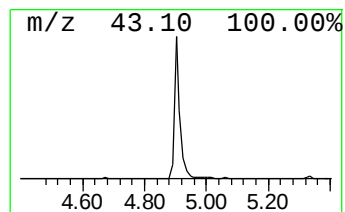
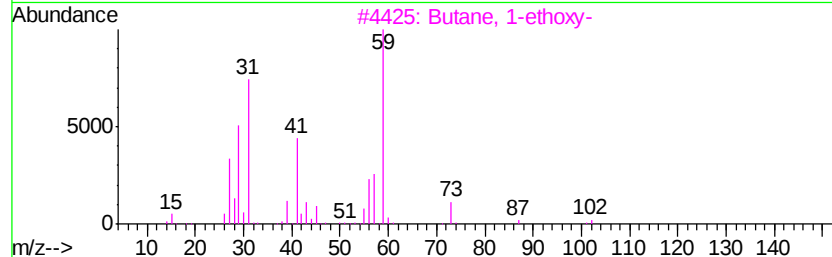
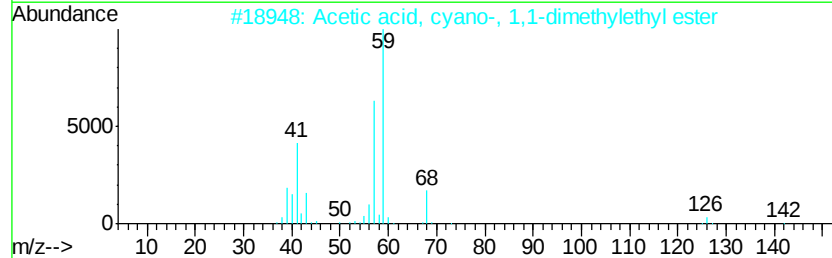
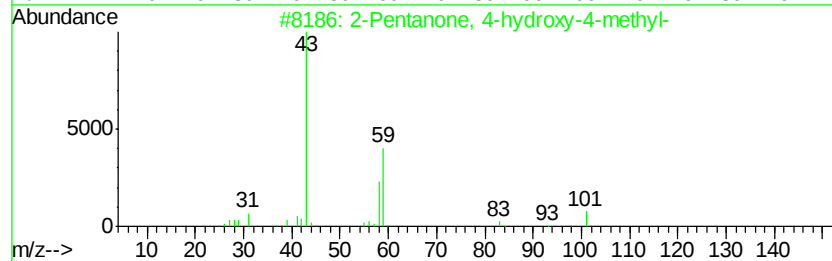
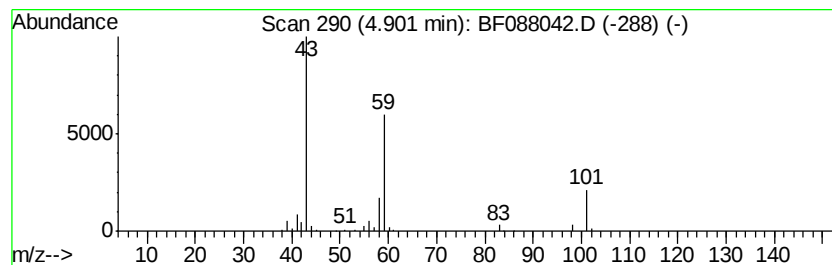
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Peak Number 3 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.90	6.56 ng	195092	1,4-Dichlorobenzene-d4	6.74

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56
2		Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	17
3		Butane, 1-ethoxy-	102	C6H14O	000628-81-9	9
4		Acetone	58	C3H6O	000067-64-1	9
5		(+)-4-Amino-4,5-dihydro-2(3H)-f...	101	C4H7NO2	016504-58-8	9



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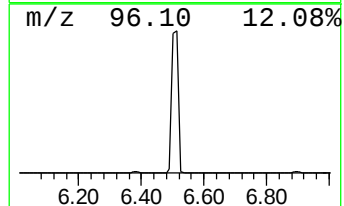
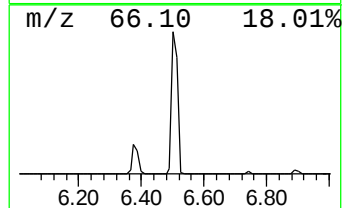
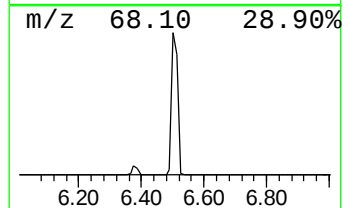
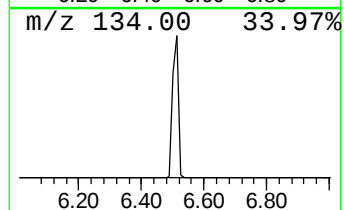
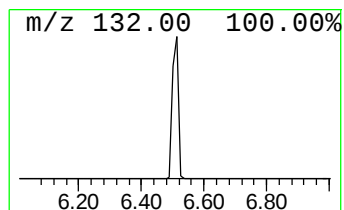
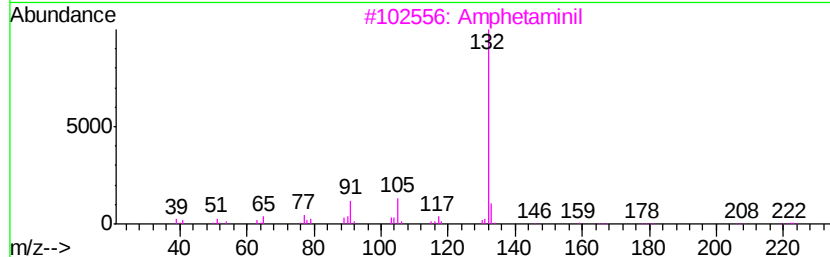
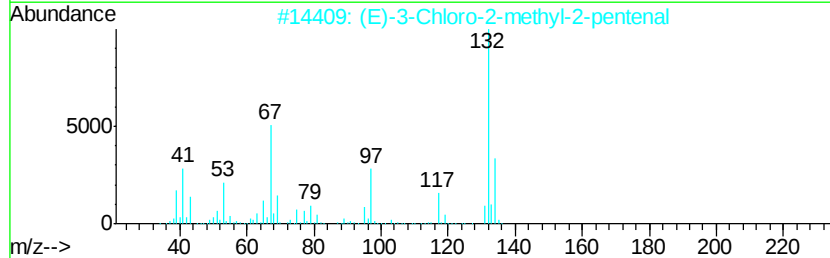
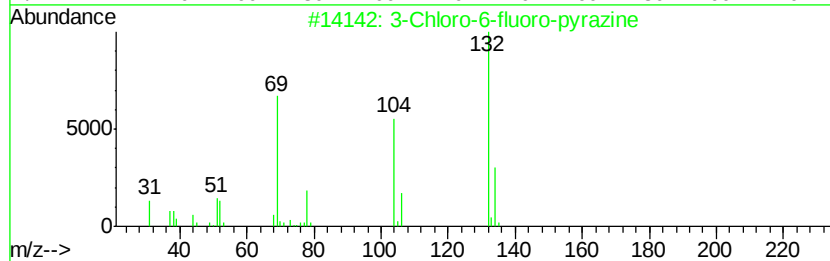
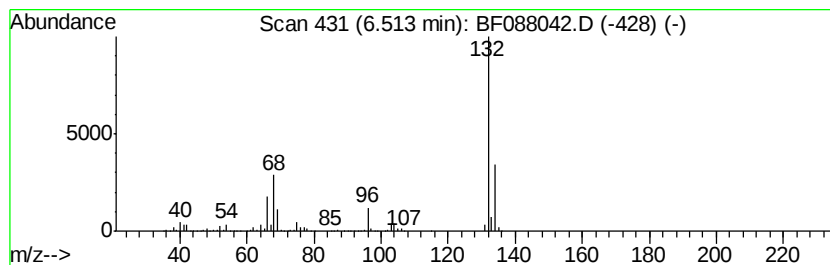
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Peak Number 4 unknown6.51 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.51	55.98 ng	1666030	1,4-Dichlorobenzene-d4	6.74

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	40
3		Amphetaminil	250	C17H18N2	017590-01-1	28
4		1H-Indazole, 7-methyl-	132	C8H8N2	1000316-00-7	25
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...	4.90	6.6	ng	195092	1	6.74	595193	20.0
unknown6.51	6.51	56.0	ng	1666030	1	6.74	595193	20.0