

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF062116\
 Data File : BF088229.D
 Acq On : 21 Jun 2016 23:10
 Operator : UM/SJ
 Sample : PB91486BL
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB91486BL

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	1.644	3	5	13	rVB	333244	508101	20.62%	2.524%
2	1.758	13	15	23	rBV	872014	1327324	53.88%	6.593%
3	4.810	280	282	291	rBV	105589	158399	6.43%	0.787%
4	5.244	318	320	323	rBV	1455718	1692641	68.70%	8.408%
5	6.307	411	413	415	rBV	1713509	1807891	73.38%	8.981%
6	6.422	421	423	425	rBV	1970595	1851680	75.16%	9.198%
7	6.650	441	443	446	rBV	576133	474332	19.25%	2.356%
8	6.810	454	457	459	rBV	1199333	1587949	64.46%	7.888%
9	7.222	491	493	495	rBV	1265290	1270677	51.58%	6.312%
10	7.930	553	555	558	rBV	616009	637360	25.87%	3.166%
11	9.016	647	650	652	rBV	1944612	2463655	100.00%	12.238%
12	9.679	706	708	711	rBV	681450	724208	29.40%	3.597%
13	10.468	775	777	780	rBV2	1048253	1304661	52.96%	6.481%
14	11.165	836	838	840	rBV	721667	742275	30.13%	3.687%
15	12.742	974	976	979	rBV	1749673	2348856	95.34%	11.668%
16	13.794	1065	1068	1070	rBV	578171	704470	28.59%	3.499%
17	15.165	1185	1188	1196	rBV	410621	526600	21.37%	2.616%

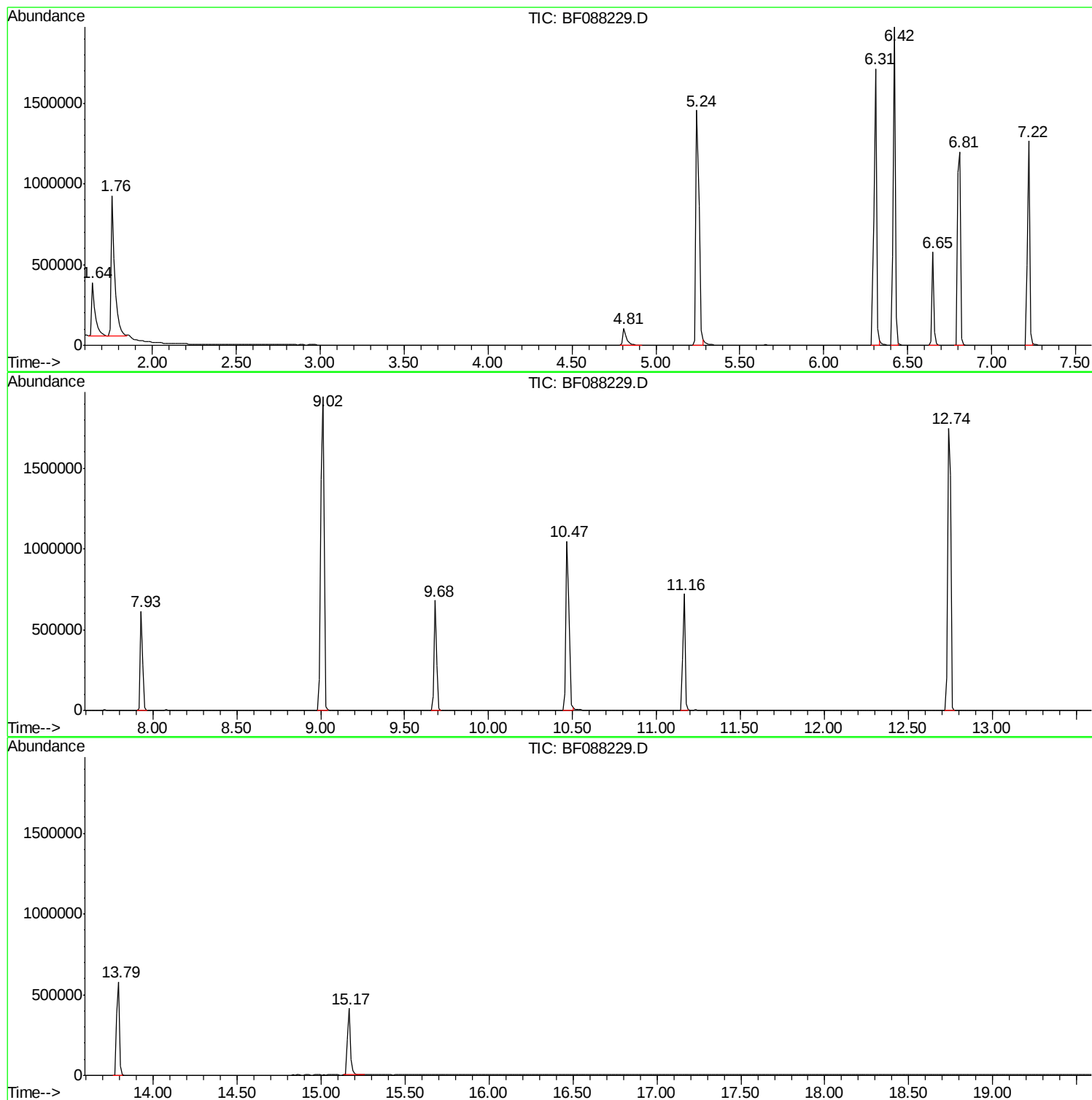
Sum of corrected areas: 20131079

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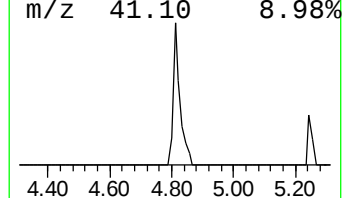
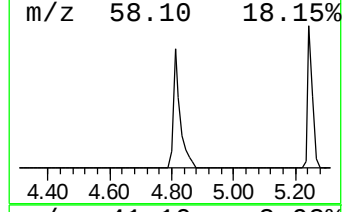
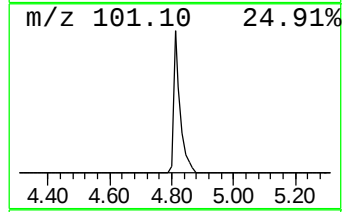
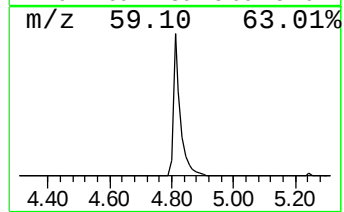
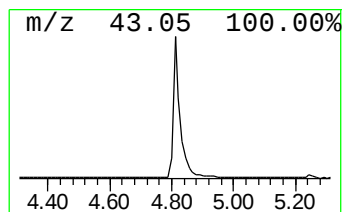
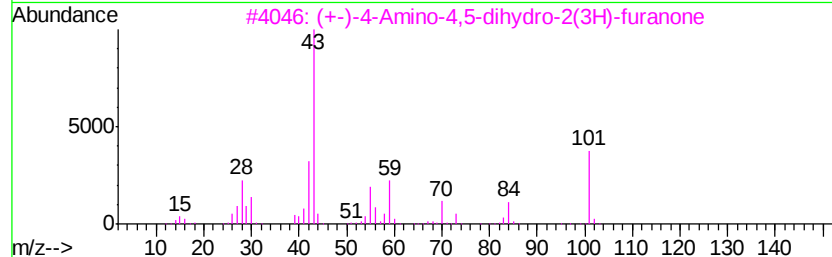
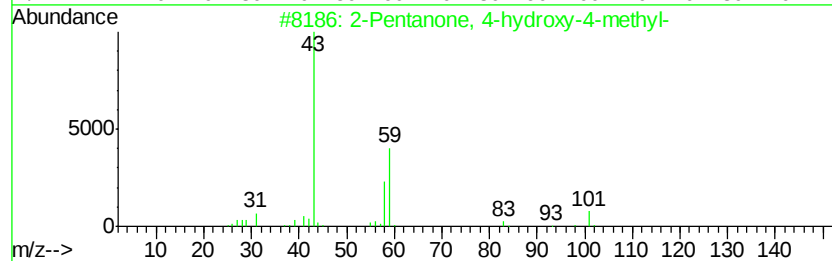
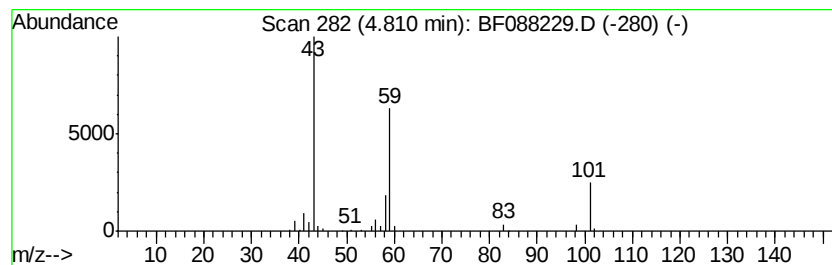
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Peak Number 3 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.81	6.68 ng	158399	1,4-Dichlorobenzene-d4	6.65

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	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	(+)-4-Amino-4,5-dihydro-2(3H)-f...	101	C4H7NO2		016504-58-8	9
4	5-Hexen-2-one	98	C6H10O		000109-49-9	9
5	Acetone	58	C3H6O		000067-64-1	9



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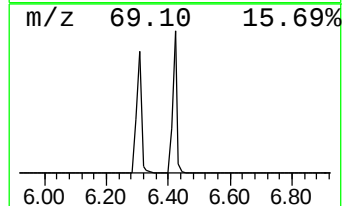
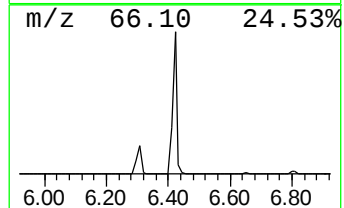
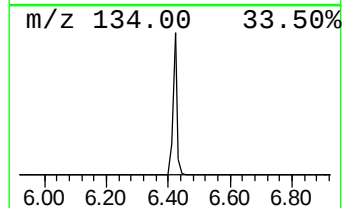
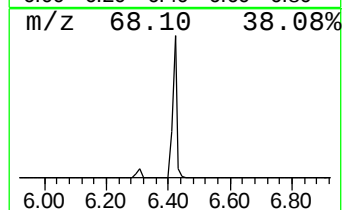
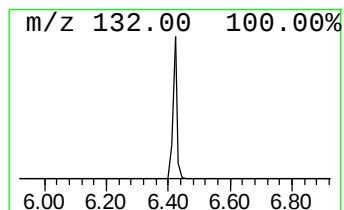
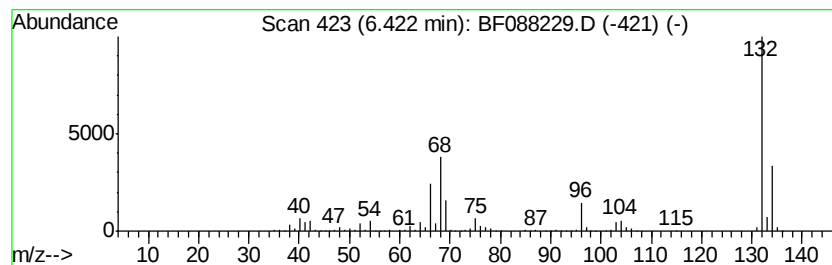
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Peak Number 4 unknown6.42 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.42	78.08 ng	1851680	1,4-Dichlorobenzene-d4	6.65

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Chloro-6-fluoro-pyrazine			132	C4H2ClFN2	1000146-10-7	27
2	(E)-3-Chloro-2-methyl-2-pentenal			132	C6H9ClO	031357-76-3	16
3	5-Aminoindole			132	C8H8N2	005192-03-0	12
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...	4.81	6.7	ng	158399	1	6.65	474332	20.0
unknown6.42	6.42	78.1	ng	1851680	1	6.65	474332	20.0