

Data Path : Z:\HPCHEM1\BNA G\DATA\BG050115\
 Data File : BG016845.D
 Acq On : 1 May 2015 15:19
 Operator : TP/IZ
 Sample : G2012-03
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 FYMW-06

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_G\METHODS\8270-BG042915.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	4.381	281	288	295	rBV2	28988	46087	1.40%	0.297%
2	4.863	365	370	392	rBV	308069	555206	16.88%	3.584%
3	6.479	640	645	659	rBV	171280	357016	10.86%	2.304%
4	6.802	693	700	720	rBV	734667	1240565	37.72%	8.007%
5	7.260	772	778	793	rBV	173175	317249	9.65%	2.048%
6	7.571	823	831	845	rBV	678904	1132542	34.44%	7.310%
7	8.423	969	976	994	rBV	475927	912020	27.73%	5.887%
8	10.033	1241	1250	1263	rBV	258686	476241	14.48%	3.074%
9	12.554	1671	1679	1699	rBV	1480577	2381189	72.40%	15.370%
10	13.940	1909	1915	1929	rBV2	429421	679364	20.66%	4.385%
11	15.450	2165	2172	2192	rBV	1012375	1554062	47.25%	10.031%
12	16.696	2378	2384	2397	rBV2	531049	816724	24.83%	5.272%
13	19.358	2831	2837	2848	rBV	2803622	3288738	100.00%	21.228%
14	20.656	3053	3058	3064	rBV3	35142	59156	1.80%	0.382%
15	20.909	3095	3101	3110	rBV	687949	855218	26.00%	5.520%
16	22.977	3447	3453	3465	rBV	469816	821266	24.97%	5.301%

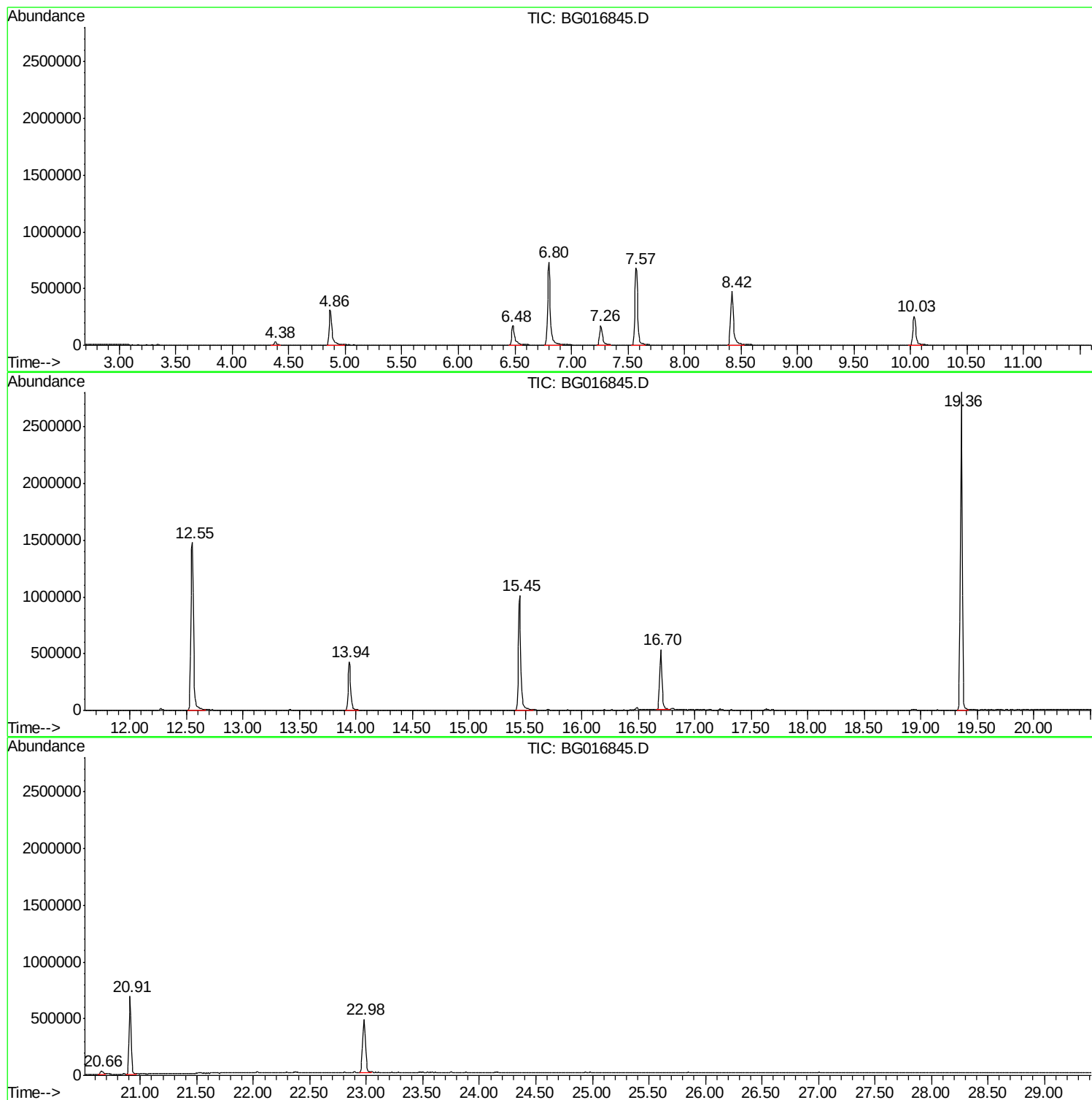
Sum of corrected areas: 15492643

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Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG042915.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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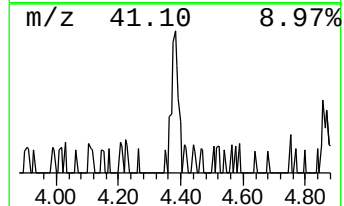
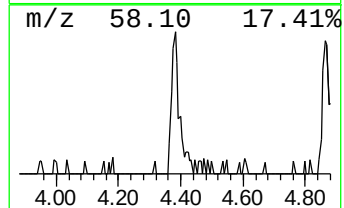
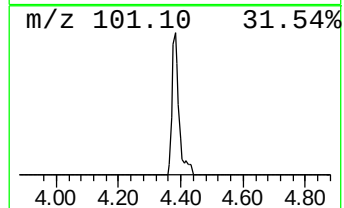
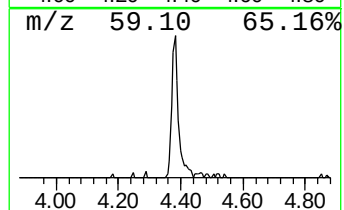
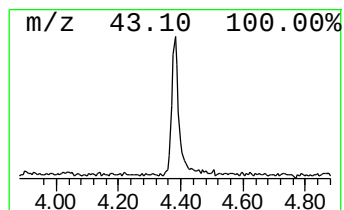
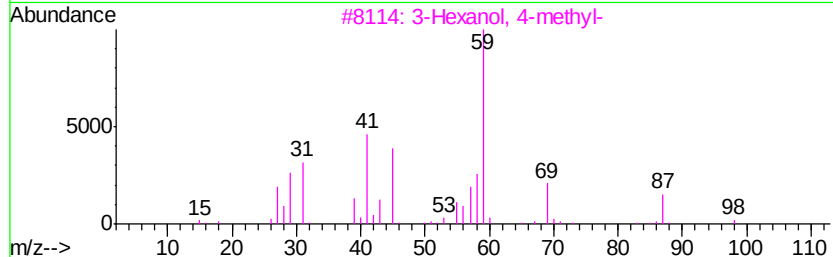
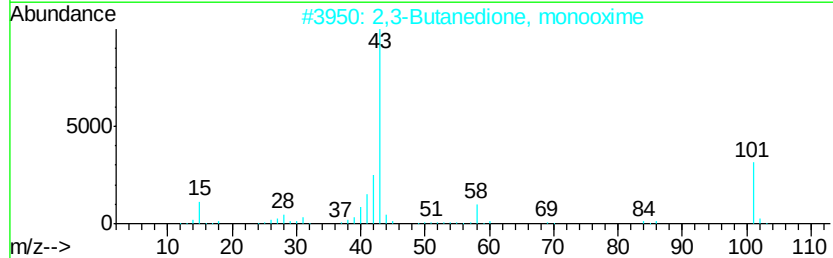
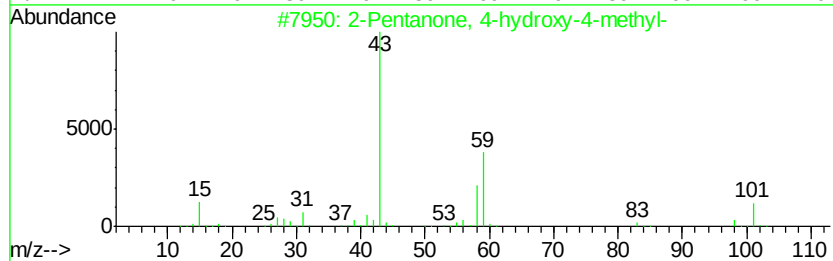
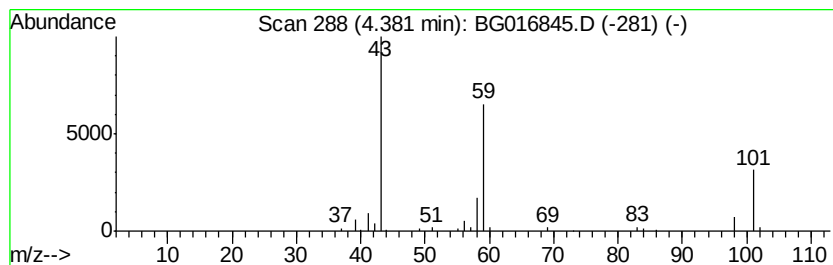
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.38	2.91 ng	46087	1,4-Dichlorobenzene-d4	7.26

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	40
2			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	35
3			3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	25
4			Butane, 1-ethoxy-	102	C6H14O	000628-81-9	9
5			(+)-4-Amino-4,5-dihydro-2(3H)-f...	101	C4H7NO2	016504-58-8	9



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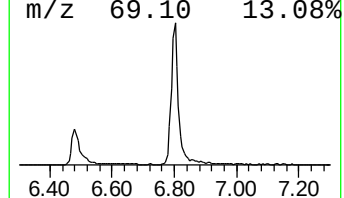
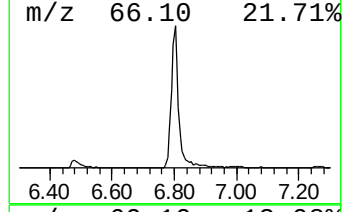
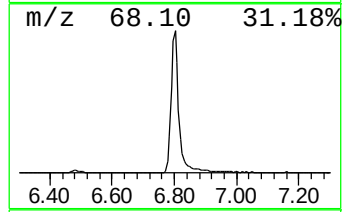
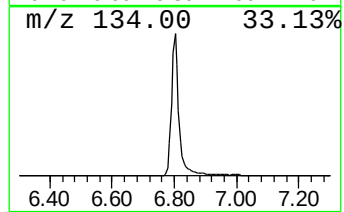
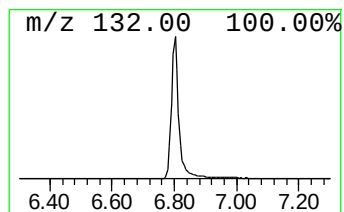
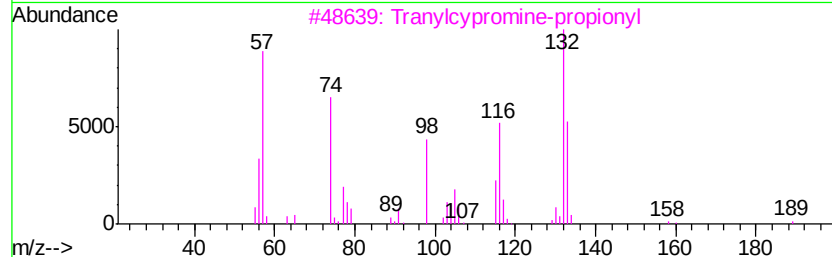
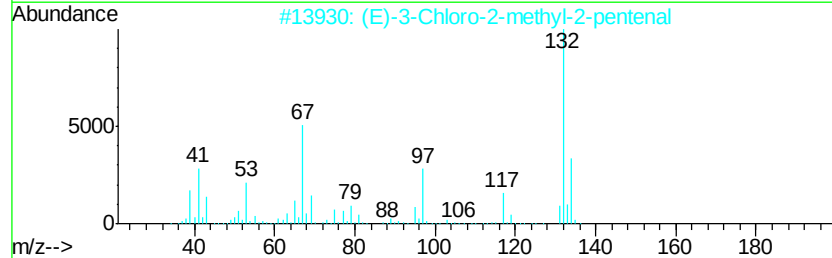
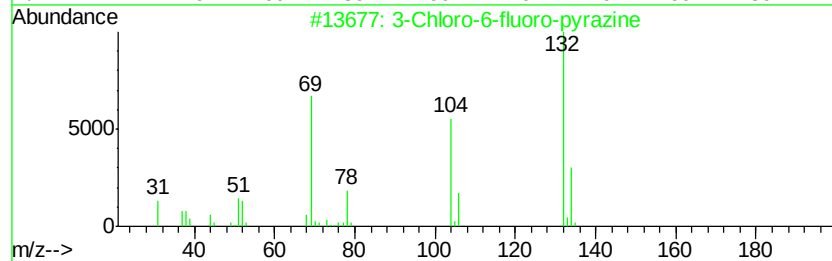
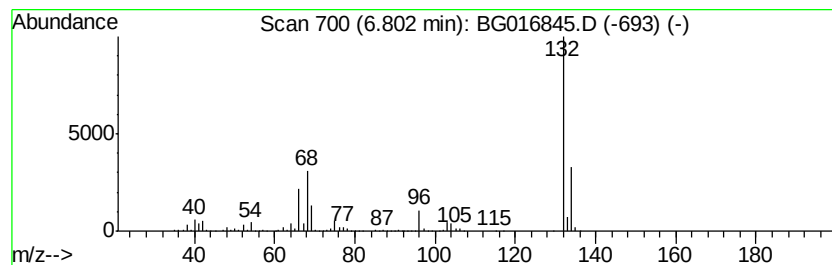
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Peak Number 2 unknown6.80 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.80	78.21 ng	1240570	1,4-Dichlorobenzene-d4	7.26

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	38
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	35
3		Tranvlcyromine-propionyl	189	C12H15NO	1000123-86-3	12
4		Xenon	132	Xe	007440-63-3	9
5		2H-Cyclopenta[d]pyridazine, 2-me...	132	C8H8N2	022291-85-6	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	4.38	2.9	ng	46087	1	7.26	317249	20.0
unknown6.80	6.80	78.2	ng	1240570	1	7.26	317249	20.0