

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA_G\DATA\BG010318\
 Data File : BG031891.D
 Acq On : 3 Jan 2018 15:22
 Operator : SJ/JU
 Sample : PB105448BL
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB105448BL

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-BG122117.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	5.697	405	410	420	rBV2	19739	35241	1.10%	0.260%
2	6.208	490	497	508	rBV	435834	747557	23.38%	5.512%
3	7.888	774	783	792	rBV	434398	779831	24.38%	5.750%
4	8.294	845	852	862	rBV	442944	826916	25.86%	6.097%
5	8.787	928	936	947	rBB	114144	215320	6.73%	1.588%
6	9.122	984	993	1005	rBB	373029	684356	21.40%	5.046%
7	9.980	1131	1139	1149	rBV	246284	457297	14.30%	3.372%
8	11.672	1419	1427	1436	rBV	163905	314563	9.84%	2.319%
9	14.046	1823	1831	1839	rBV	976156	1563818	48.90%	11.530%
10	15.415	2057	2064	2074	rVB2	303811	516609	16.15%	3.809%
11	16.889	2307	2315	2325	rBV	861410	1368432	42.79%	10.090%
12	18.147	2522	2529	2539	rBV2	497930	791182	24.74%	5.833%
13	20.679	2954	2960	2971	rBV	2416195	3198053	100.00%	23.579%
14	22.489	3258	3268	3276	rBV2	549964	1032795	32.29%	7.615%
15	26.220	3890	3903	3916	rBV2	320313	1030946	32.24%	7.601%

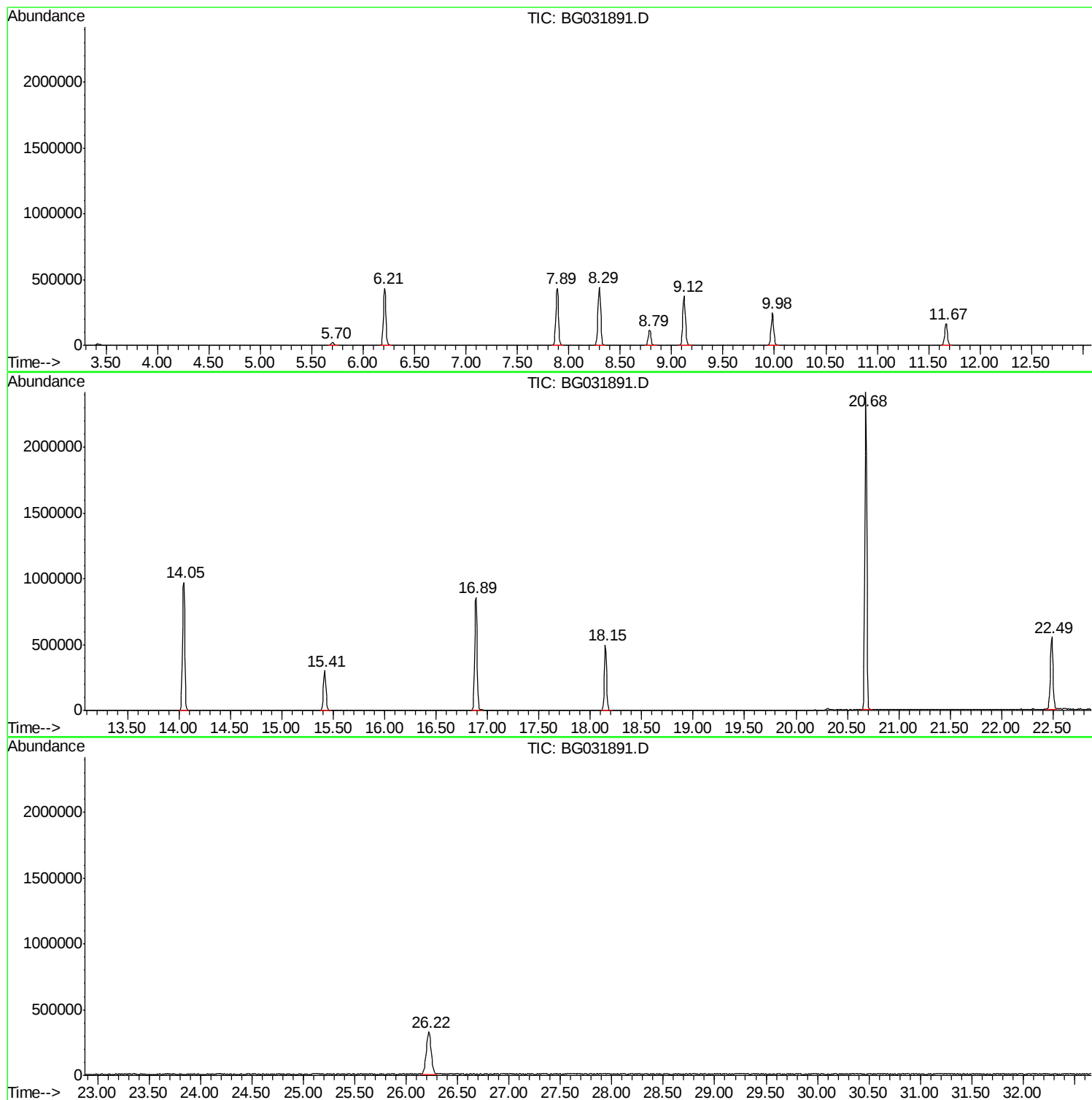
Sum of corrected areas: 13562916

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Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG122117.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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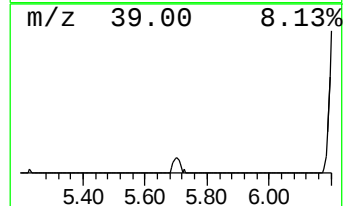
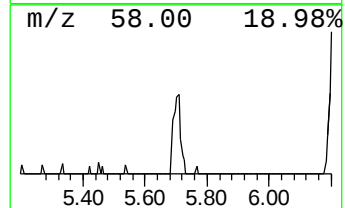
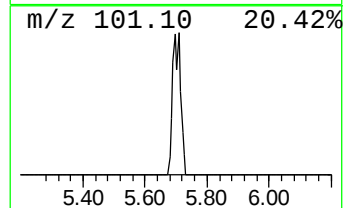
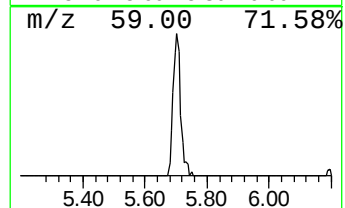
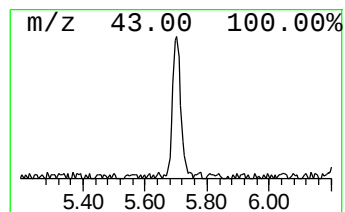
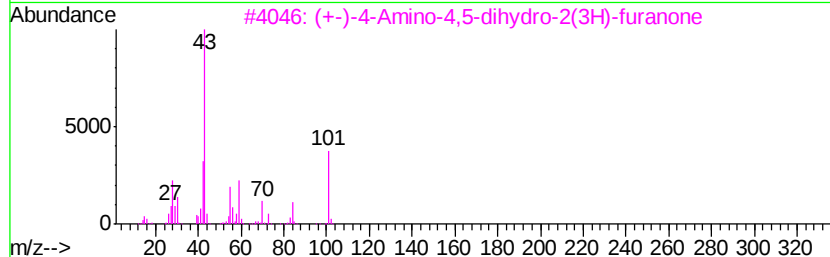
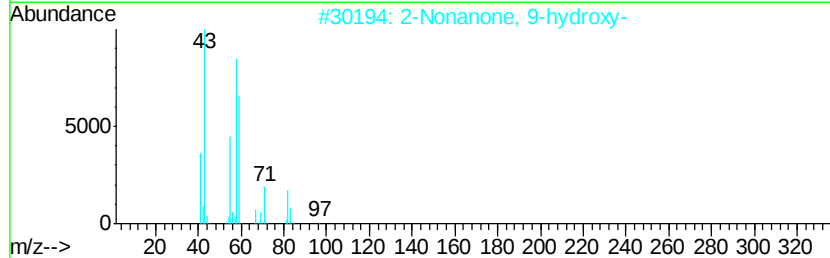
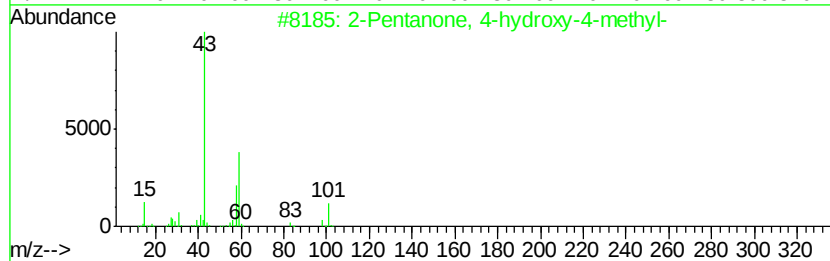
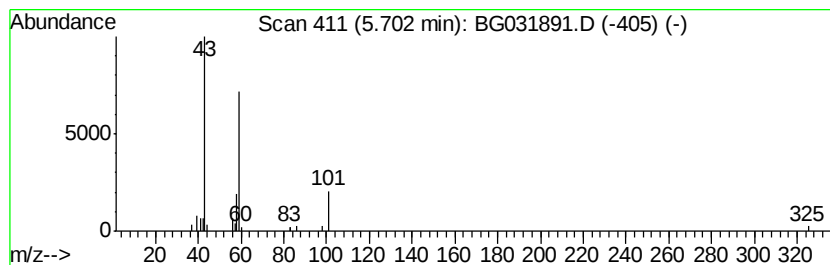
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.70	3.27 ng	35241	1,4-Dichlorobenzene-d4	8.79

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	40
2			2-Nonanone, 9-hydroxy-	158	C9H18O2	025368-56-3	28
3			(+)-4-Amino-4,5-dihydro-2(3H)-f...	101	C4H7NO2	016504-58-8	23
4			4-Penten-2-one, 3-methyl-	98	C6H10O	000758-87-2	9
5			5-Hexen-2-one	98	C6H10O	000109-49-9	9



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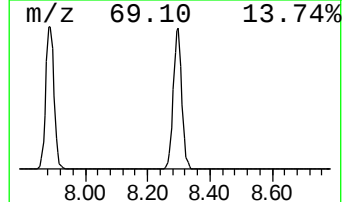
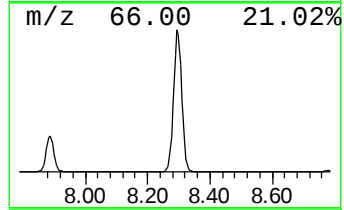
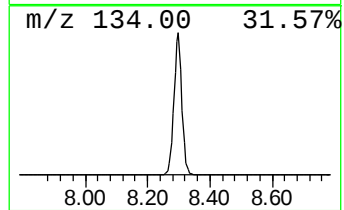
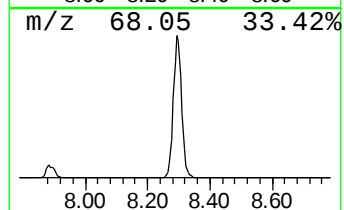
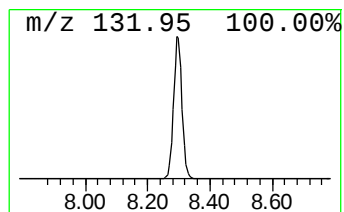
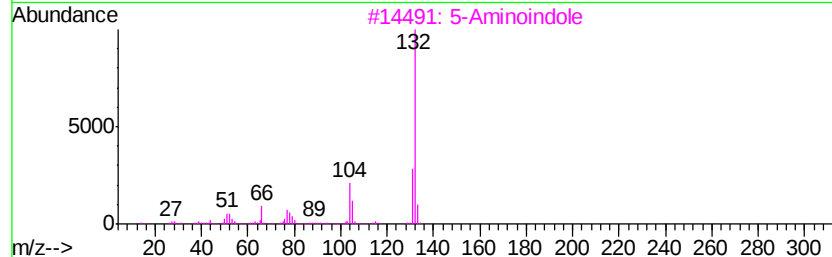
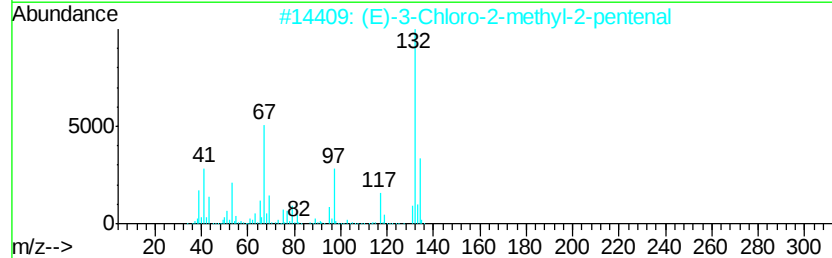
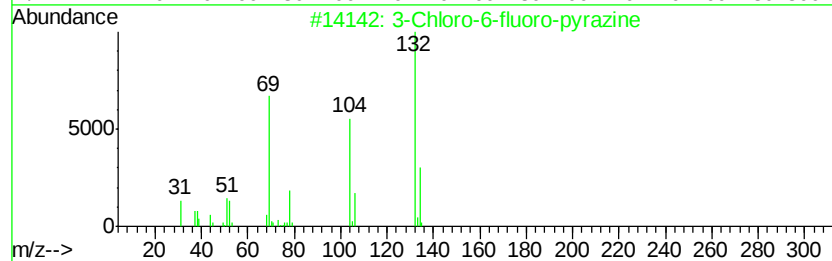
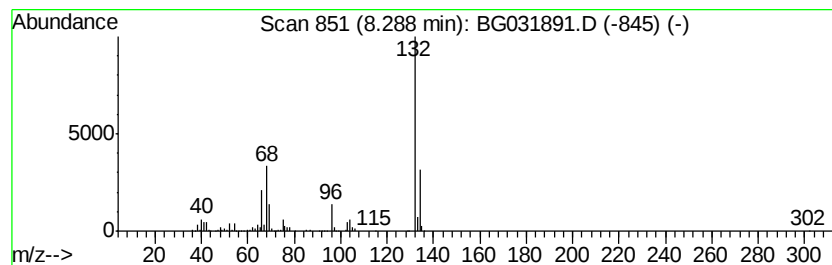
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Peak Number 2 unknown8.29 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.29	76.81 ng	826916	1,4-Dichlorobenzene-d4	8.79

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		5-Aminoindole	132	C8H8N2	005192-03-0	12
4		1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	10
5		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.70	3.3	ng	35241	1	8.79	215320	20.0
unknown8.29	8.29	76.8	ng	826916	1	8.79	215320	20.0