

Data Path : Z:\HPCHEM1\BNA G\DATA\BG042018\
 Data File : BG033900.D
 Acq On : 21 Apr 2018 1:19
 Operator : SJ/JU
 Sample : J2473-03
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 041818-DBRI-F-U-B

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-BG041918.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	4.995	283	290	297	rBV	66656	105416	1.64%	0.448%
2	5.442	358	366	378	rBV	463815	734031	11.40%	3.120%
3	6.993	621	630	638	rBV	294828	494828	7.68%	2.103%
4	7.363	686	693	705	rVB	801945	1343250	20.86%	5.709%
5	7.833	765	773	780	rBV	231975	402343	6.25%	1.710%
6	8.150	816	827	836	rVB	996316	1691897	26.27%	7.191%
7	8.979	958	968	978	rBV	850852	1483554	23.04%	6.306%
8	10.612	1239	1246	1255	rVB	329753	588290	9.14%	2.500%
9	13.086	1660	1667	1676	rBV	2249431	3509350	54.50%	14.916%
10	14.467	1893	1902	1913	rBV2	512607	863216	13.40%	3.669%
11	15.953	2148	2155	2166	rVB	1269010	1891847	29.38%	8.041%
12	17.216	2363	2370	2376	rBV	739046	1075114	16.70%	4.570%
13	18.133	2520	2526	2536	rBV3	48172	85928	1.33%	0.365%
14	19.854	2812	2819	2825	rBV	4804563	6439549	100.00%	27.371%
15	21.470	3087	3094	3104	rVB2	906948	1435725	22.30%	6.102%
16	24.537	3605	3616	3629	rVB2	520406	1382709	21.47%	5.877%

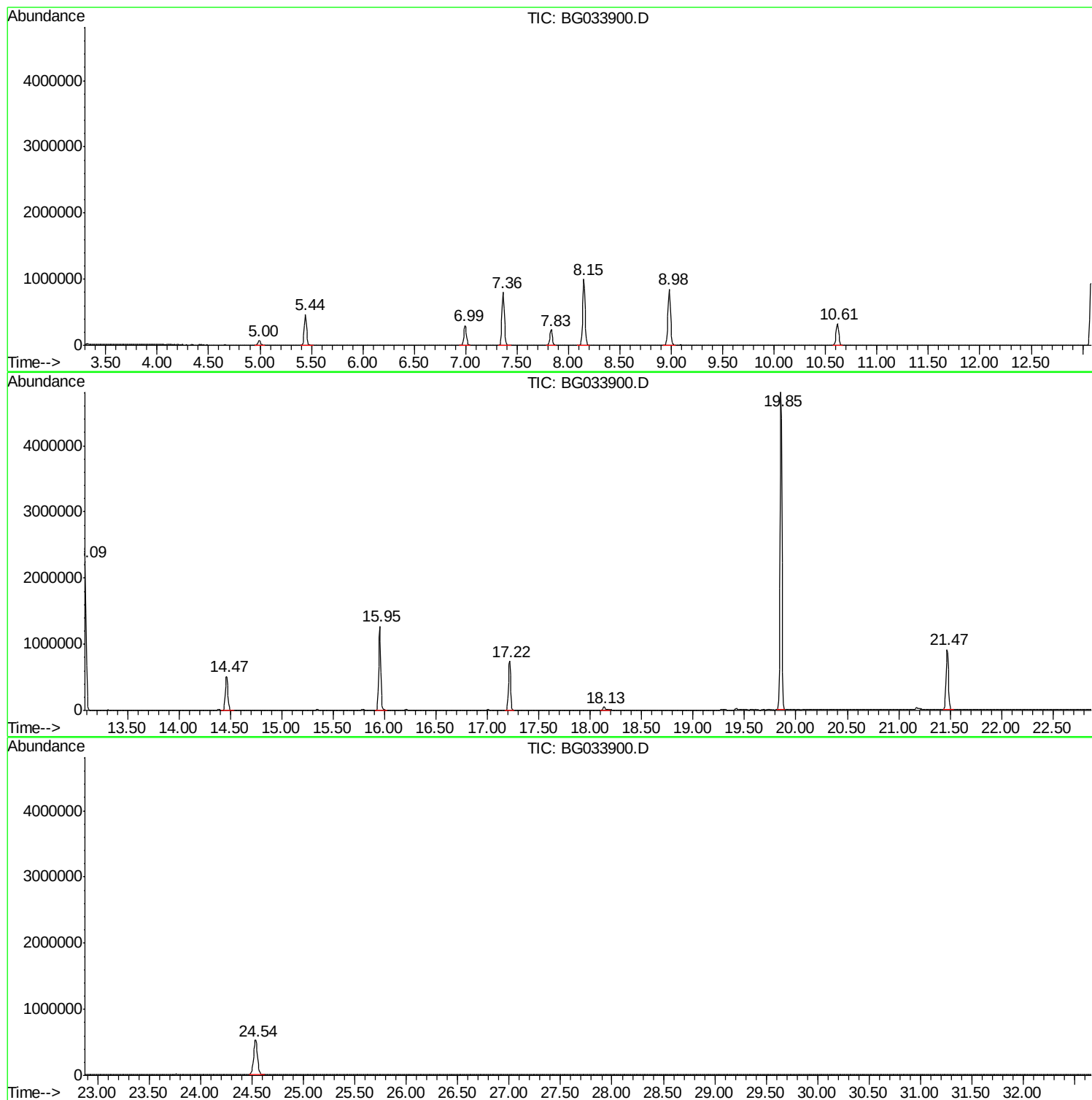
Sum of corrected areas: 23527047

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Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG041918.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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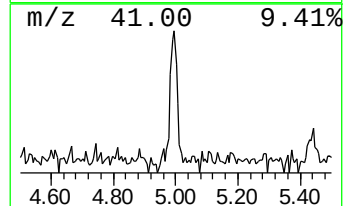
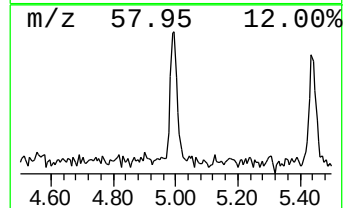
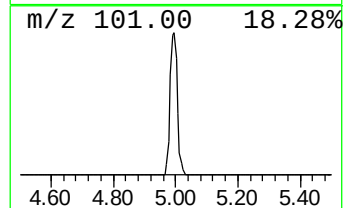
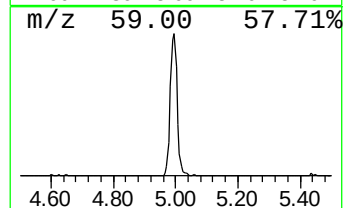
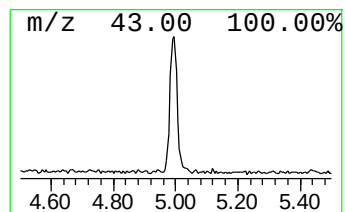
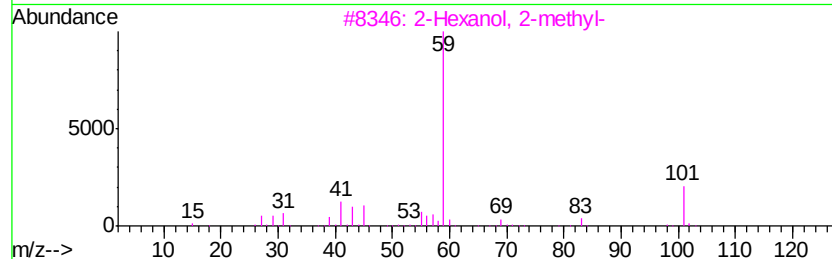
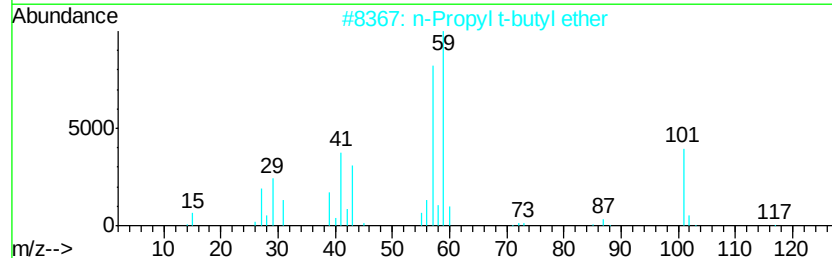
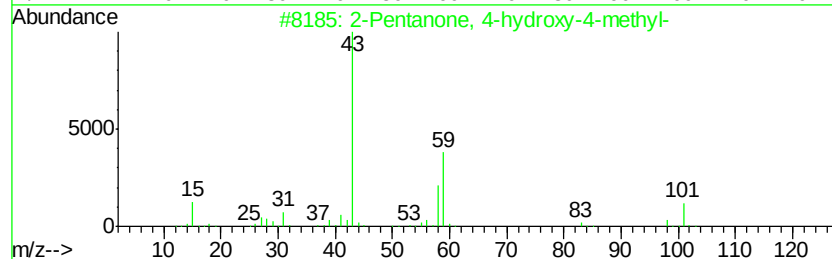
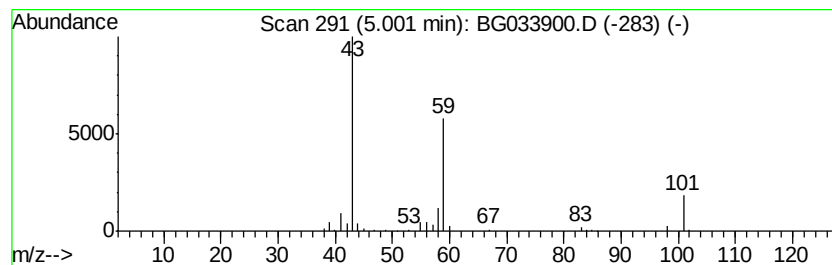
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.00	5.24 ng	105416	1,4-Dichlorobenzene-d4	7.83

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50
2		n-Propyl t-butyl ether	116	C7H16O	1000334-81-9	28
3		2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	28
4		2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	27
5		1,2-Butanediol	90	C4H10O2	000584-03-2	25



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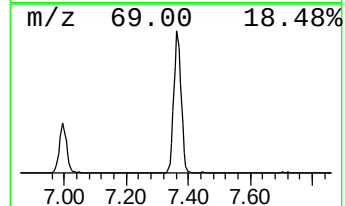
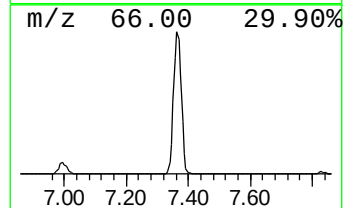
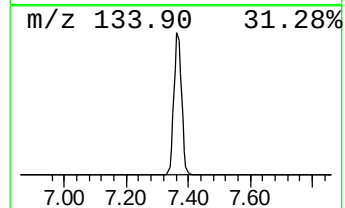
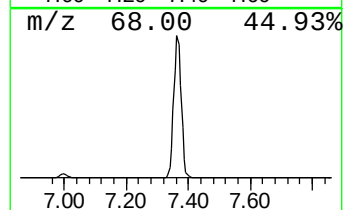
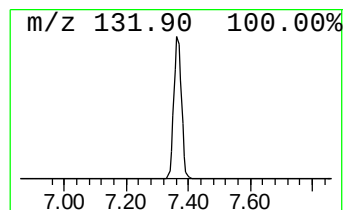
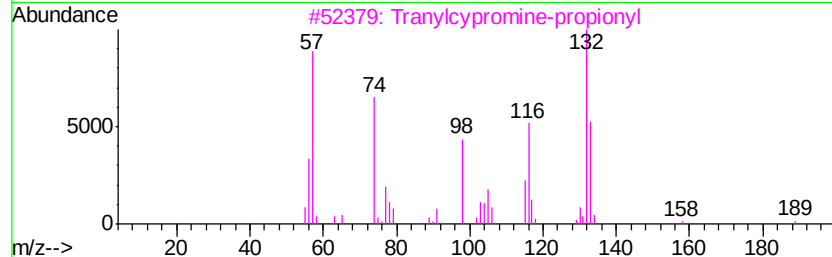
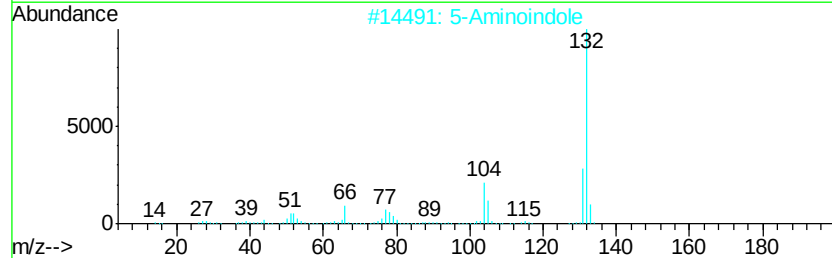
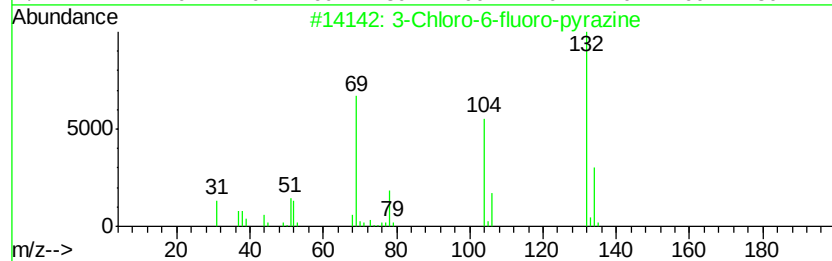
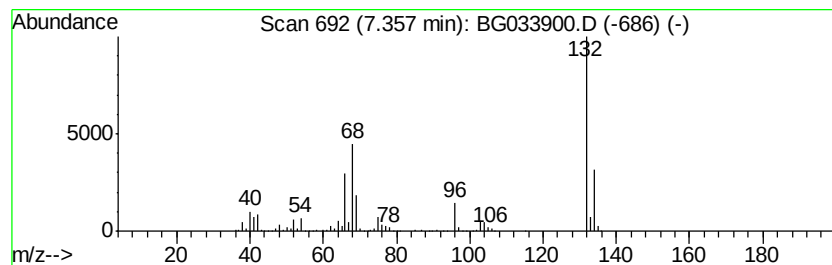
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Peak Number 2 unknown7.36 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.36	66.77 ng	1343250	1,4-Dichlorobenzene-d4	7.83

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Chloro-6-fluoro-pyrazine			132	C4H2ClFN2	1000146-10-7	27
2	5-Aminoindole			132	C8H8N2	005192-03-0	12
3	Tranvlcvpromine-propionyl			189	C12H15NO	1000123-86-3	10
4	3H-1,2-Dithiol-3-one, 5-methyl-			132	C4H4OS2	003620-08-4	9
5	Bicyclo[4.2.0]octa-1,3,5-triene, ...			132	C10H12	028749-81-7	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.00	5.2	ng	105416	1	7.83	402343	20.0
unknown7.36	7.36	66.8	ng	1343250	1	7.83	402343	20.0