

Data Path : Z:\HPCHEM1\BNA G\DATA\BG042018\
 Data File : BG033904.D
 Acq On : 21 Apr 2018 3:50
 Operator : SJ/JU
 Sample : J2473-07
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 041818-SIDI-F-D-S

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	4.991	284	290	297	rBV	65806	96576	1.52%	0.438%
2	5.438	359	366	374	rBV	431643	669653	10.50%	3.035%
3	6.995	622	631	640	rBV	271964	459410	7.21%	2.082%
4	7.365	686	694	707	rVB	736985	1226014	19.23%	5.556%
5	7.829	766	773	781	rVB	209452	357635	5.61%	1.621%
6	8.152	820	828	837	rVB	894318	1551574	24.34%	7.032%
7	8.981	960	969	977	rBV	782593	1344465	21.09%	6.093%
8	10.614	1239	1247	1255	rBV	289836	514414	8.07%	2.331%
9	13.088	1661	1668	1675	rBV	2107754	3222321	50.55%	14.604%
10	14.462	1895	1902	1911	rBV2	463419	748222	11.74%	3.391%
11	15.955	2149	2156	2166	rVB	1182927	1778183	27.89%	8.059%
12	17.212	2364	2370	2377	rBV	649208	1006648	15.79%	4.562%
13	19.856	2813	2820	2825	rBV	4848485	6374643	100.00%	28.890%
14	21.472	3089	3095	3103	rVB	915794	1380231	21.65%	6.255%
15	24.533	3605	3616	3629	rVB	499957	1335245	20.95%	6.051%

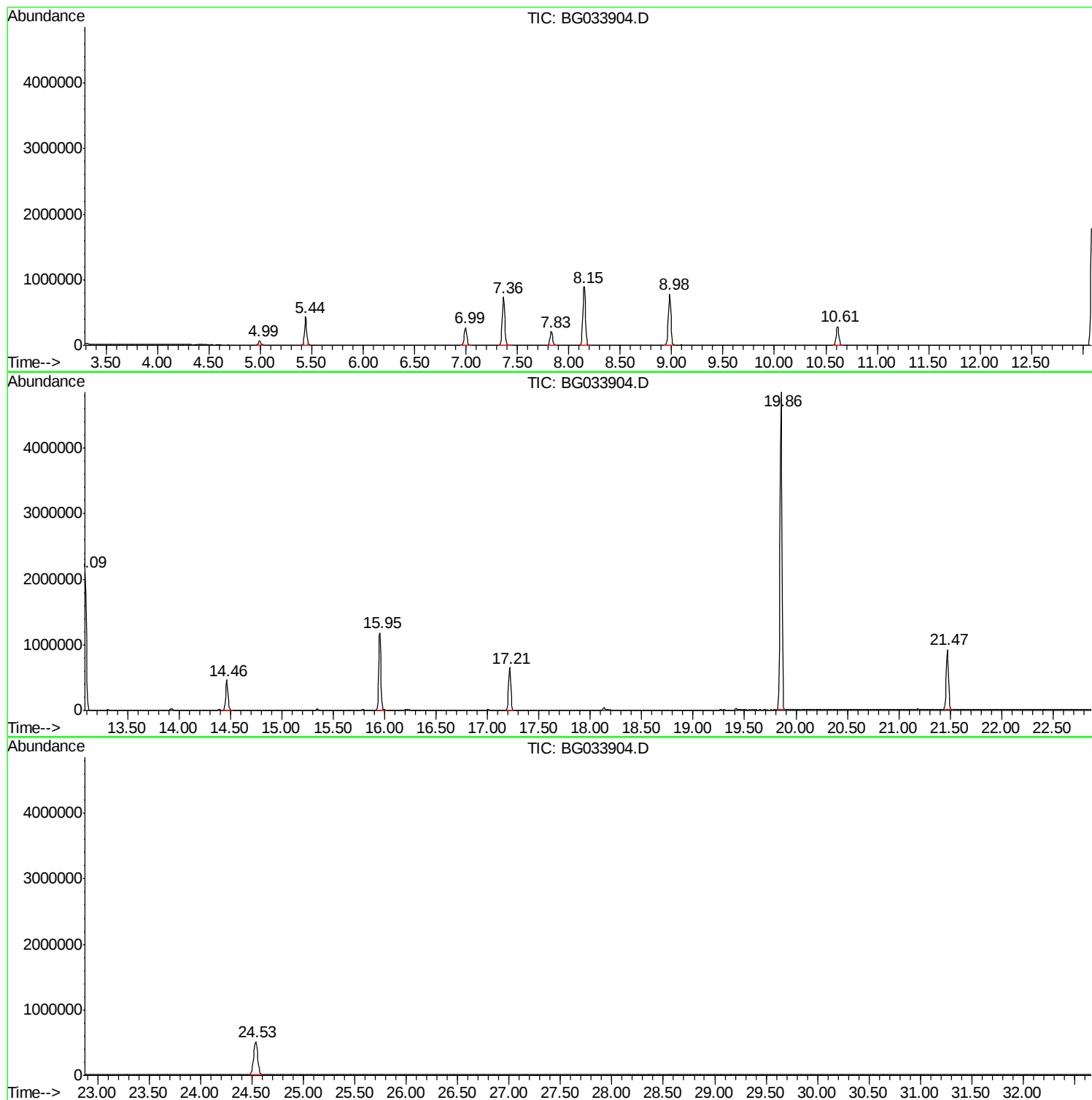
Sum of corrected areas: 22065234

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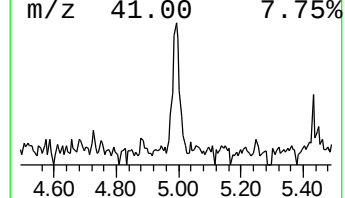
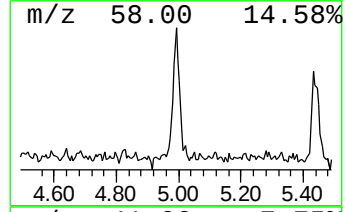
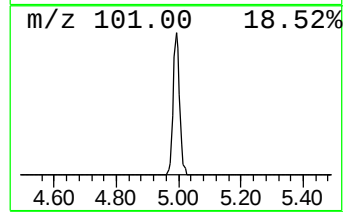
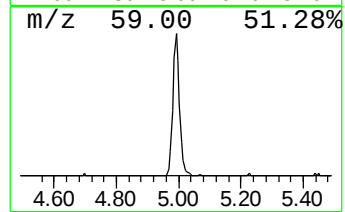
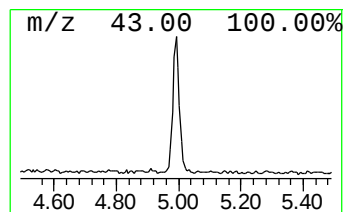
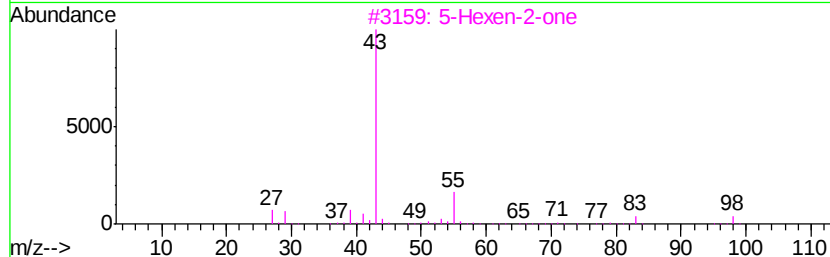
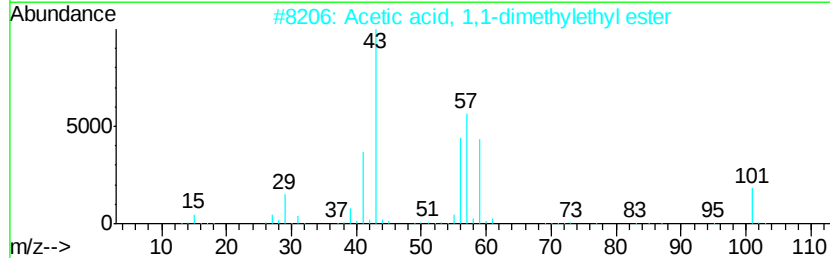
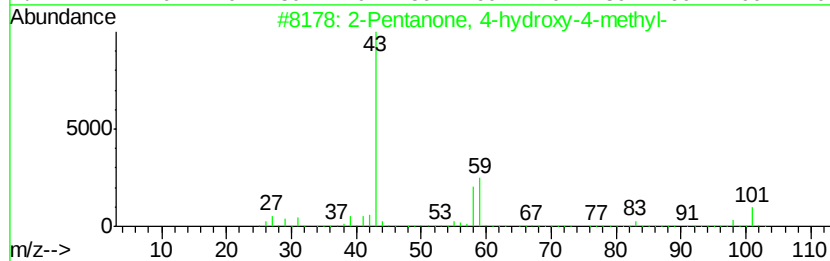
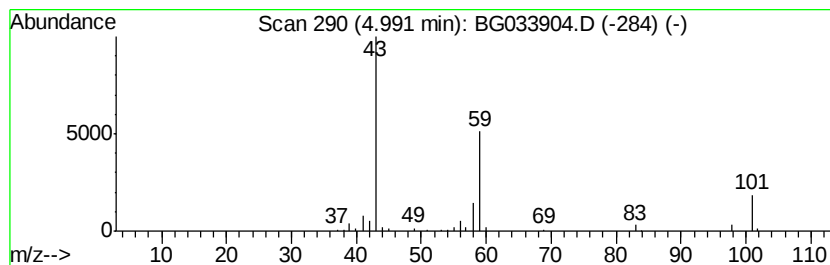
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.99	5.40 ng	96576	1,4-Dichlorobenzene-d4	7.83

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	64
2			Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	38
3			5-Hexen-2-one	98	C6H10O	000109-49-9	9
4			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9
5			Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9



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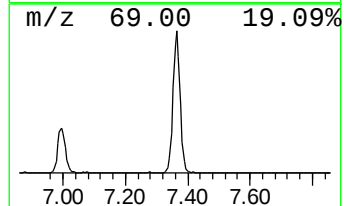
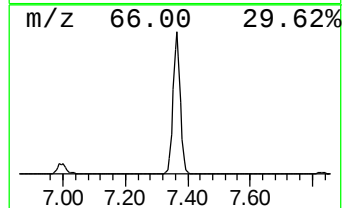
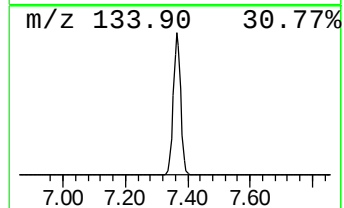
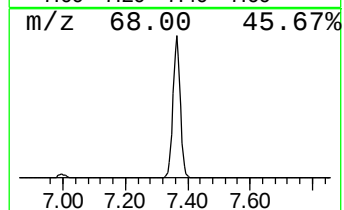
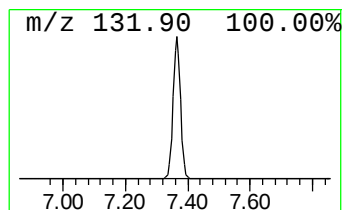
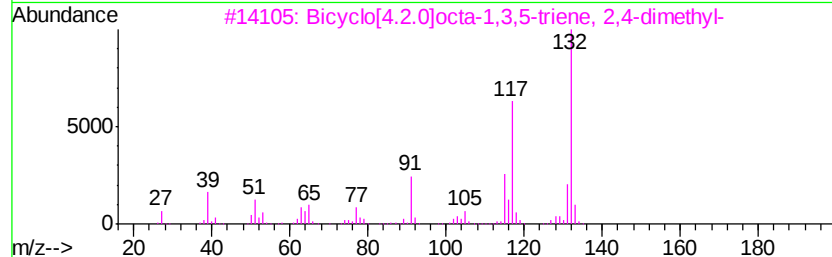
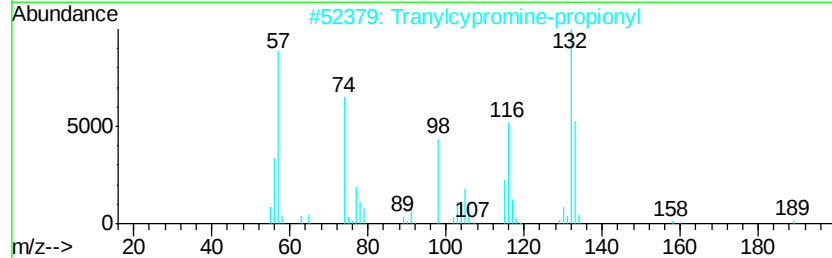
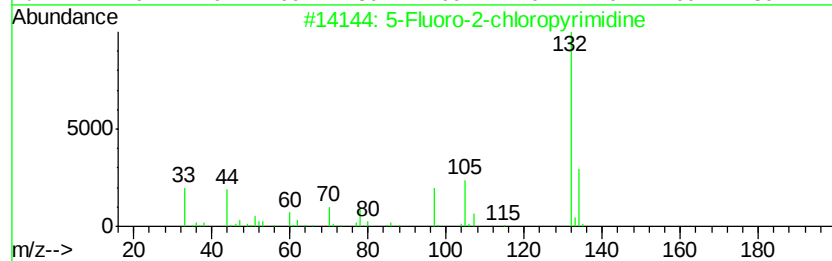
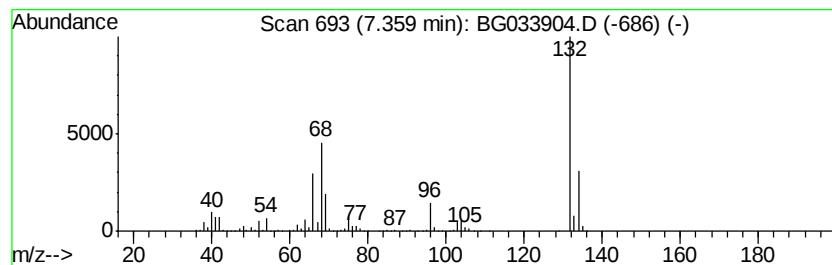
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Peak Number 2 unknown7.36 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.36	68.56 ng	1226010	1,4-Dichlorobenzene-d4	7.83

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	12
2		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
3		Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9
4		Benzene, 1-ethenyl-3,5-dimethyl-	132	C10H12	005379-20-4	9
5		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	4.99	5.4	ng	96576	1	7.83	357635	20.0
unknown7.36	7.36	68.6	ng	1226010	1	7.83	357635	20.0