

Data Path : Z:\HPCHEM1\BNA G\DATA\BG042018\
 Data File : BG033903.D
 Acq On : 21 Apr 2018 3:12
 Operator : SJ/JU
 Sample : J2473-06
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

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

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	4.992	285	290	298	rBV	70622	102318	1.51%	0.449%
2	5.438	360	366	375	rBV	426270	650570	9.63%	2.853%
3	6.995	622	631	640	rBV	269113	448767	6.64%	1.968%
4	7.365	687	694	702	rBV	720524	1209694	17.91%	5.304%
5	7.829	766	773	781	rVB	215734	362405	5.36%	1.589%
6	8.147	820	827	835	rVB	909840	1576027	23.33%	6.910%
7	8.981	961	969	979	rBV	783090	1349649	19.98%	5.918%
8	10.614	1239	1247	1257	rVB	299922	520501	7.70%	2.282%
9	13.088	1661	1668	1675	rVB	2091100	3251397	48.13%	14.256%
10	14.463	1894	1902	1911	rBV2	484117	784201	11.61%	3.438%
11	15.955	2149	2156	2163	rBV	1197989	1840735	27.25%	8.071%
12	17.212	2364	2370	2379	rVB2	713225	1069947	15.84%	4.691%
13	19.856	2813	2820	2828	rBV	4928733	6756042	100.00%	29.623%
14	21.472	3088	3095	3103	rBV2	947620	1454256	21.53%	6.376%
15	24.533	3605	3616	3631	rVB2	522727	1430385	21.17%	6.272%

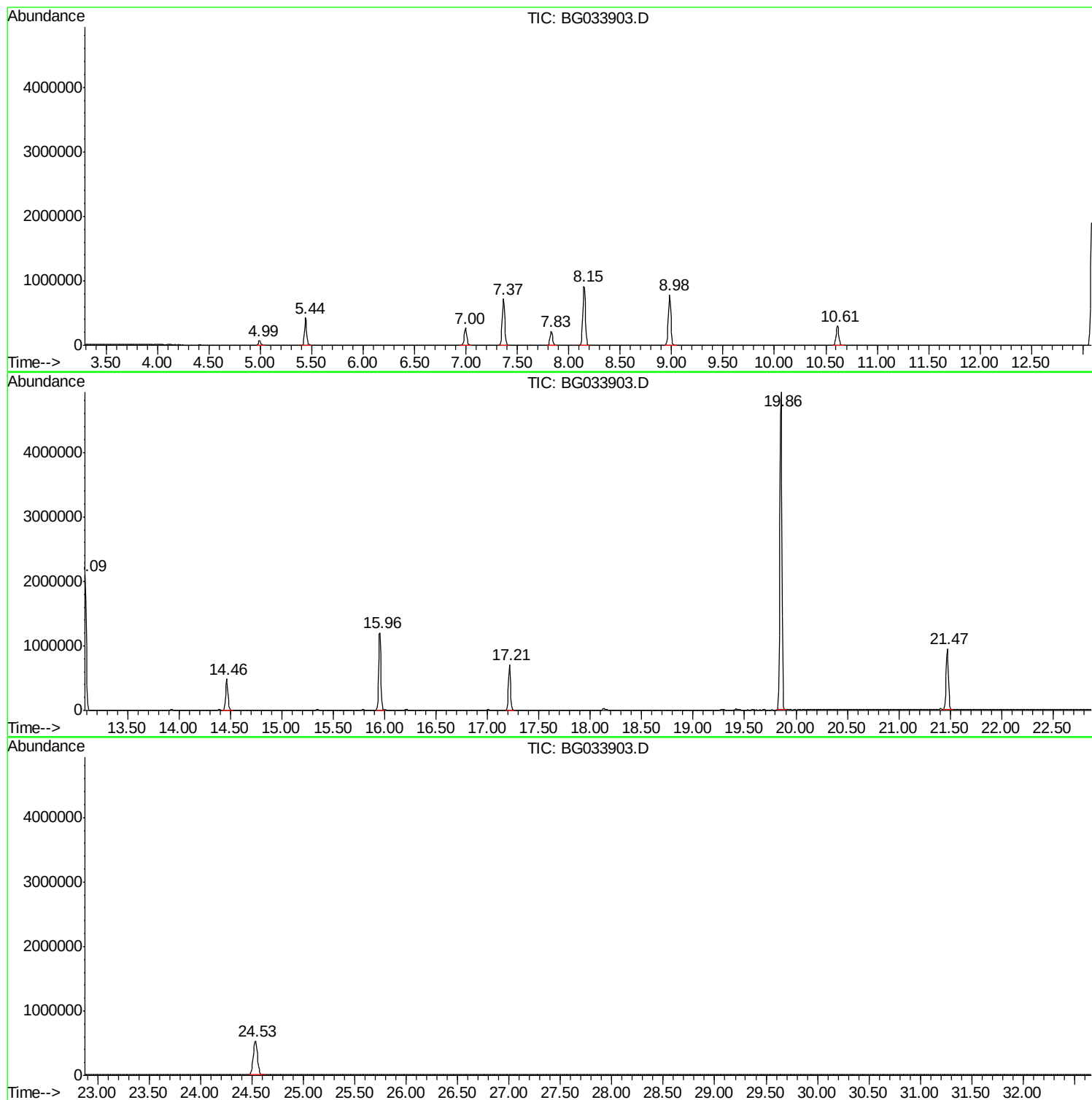
Sum of corrected areas: 22806894

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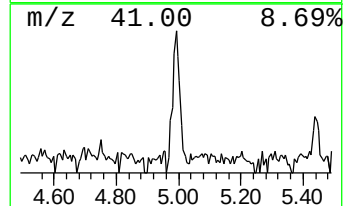
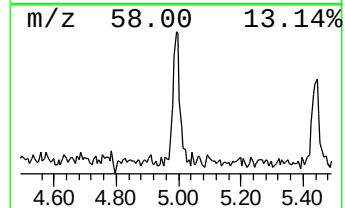
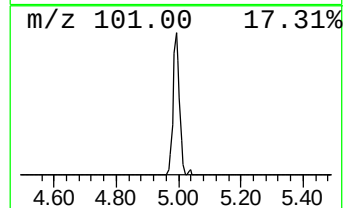
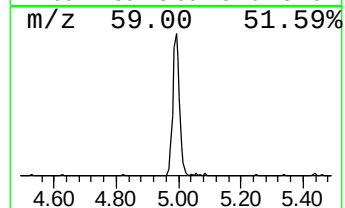
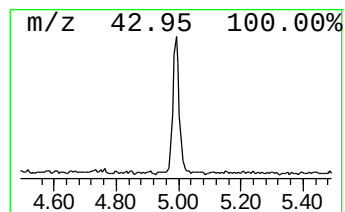
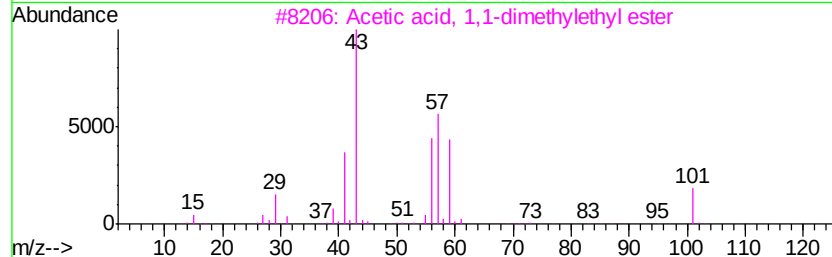
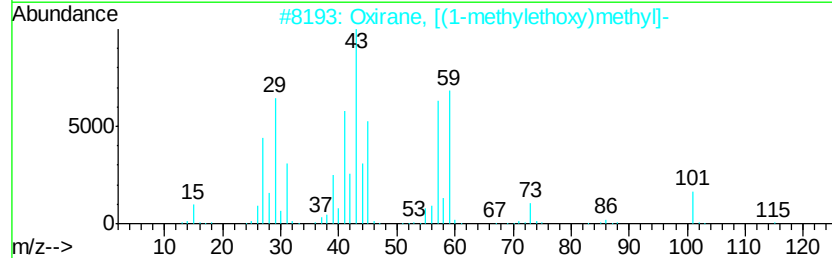
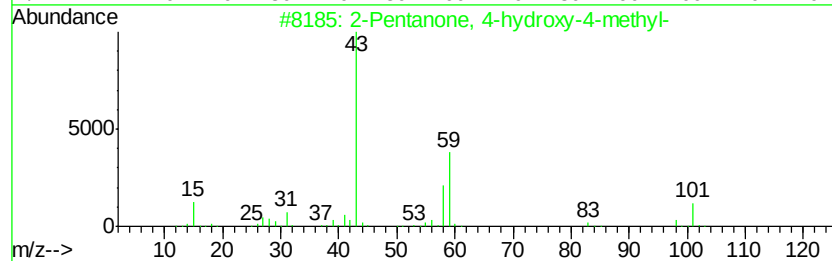
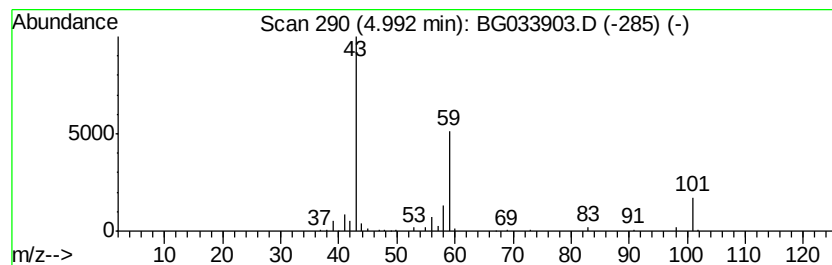
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.99	5.65 ng	102318	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	45
2			Oxirane, [(1-methylethoxy)methyl]-	116	C6H12O2	004016-14-2	40
3			Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	38
4			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	16
5			Butane	58	C4H10	000106-97-8	9



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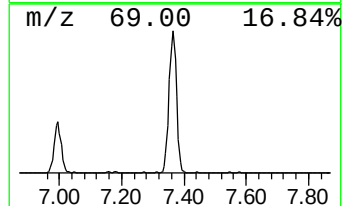
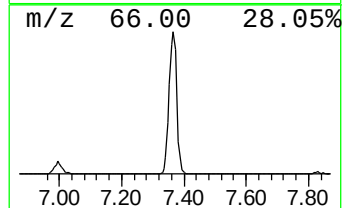
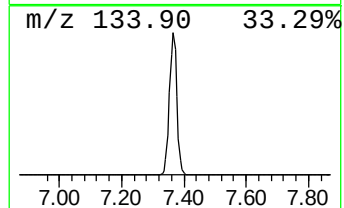
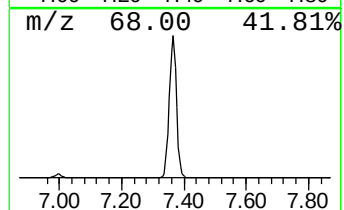
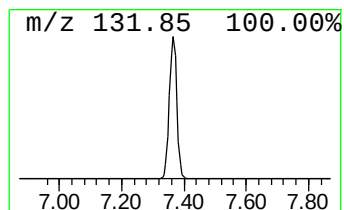
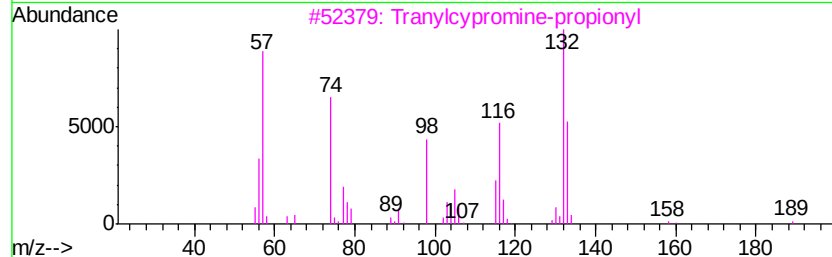
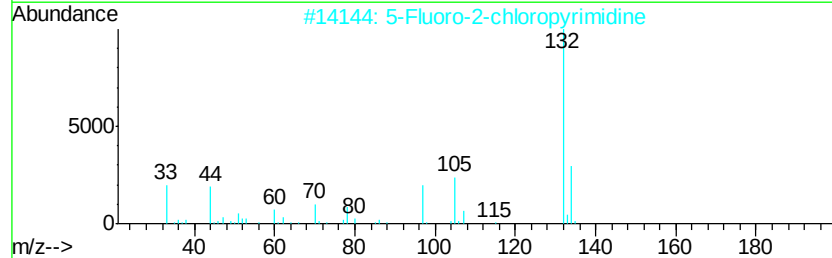
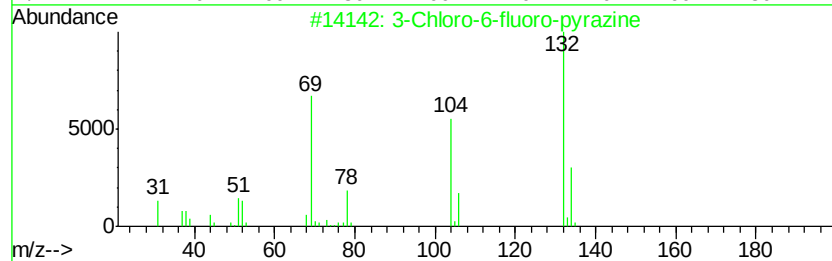
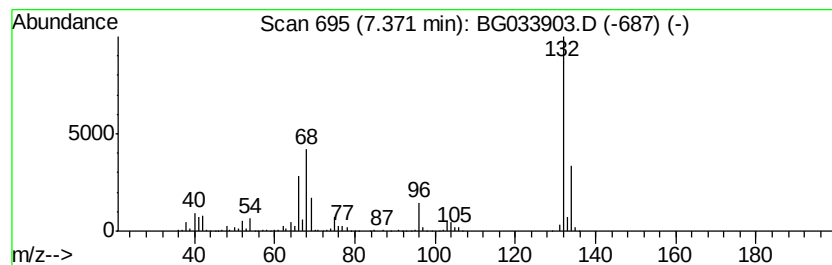
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Peak Number 2 unknown7.37 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.37	66.76 ng	1209690	1,4-Dichlorobenzene-d4	7.83

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	35
2		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	12
3		Tranvlcypropine-propionyl	189	C12H15NO	1000123-86-3	10
4		Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9
5		Benzene, 1-ethenyl-3,5-dimethyl-	132	C10H12	005379-20-4	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	4.99	5.7	ng	102318	1	7.83	362405	20.0
unknown7.37	7.37	66.8	ng	1209690	1	7.83	362405	20.0