

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110821\
 Data File : BP007875.D
 Acq On : 08 Nov 2021 23:12
 Operator : CG/JU
 Sample : M4523-07
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 FIELD-BLANK

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:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP110221.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BP007875.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	4.981	283	288	294	rBV	55516	79067	1.02%	0.288%
2	5.446	360	367	376	rBV	1129178	1602797	20.77%	5.835%
3	7.034	630	637	647	rBV	624213	987088	12.79%	3.593%
4	7.863	770	778	785	rBV	371062	587526	7.62%	2.139%
5	9.016	964	974	994	rBV	1134216	1885957	24.44%	6.866%
6	10.657	1245	1253	1261	rBV	444672	752840	9.76%	2.741%
7	13.122	1664	1672	1682	rBV	2875569	4285242	55.54%	15.600%
8	14.498	1899	1906	1916	rBV	678792	997485	12.93%	3.631%
9	15.986	2152	2159	2172	rBV	2367627	3517729	45.59%	12.806%
10	17.251	2368	2374	2387	rVB	916349	1264442	16.39%	4.603%
11	17.374	2389	2395	2400	rVB2	114729	160804	2.08%	0.585%
12	19.815	2804	2810	2815	rBV	6273265	7715348	100.00%	28.087%
13	21.339	3064	3069	3075	rBV	1314452	1743086	22.59%	6.346%
14	23.750	3473	3479	3489	rVB2	943728	1890170	24.50%	6.881%

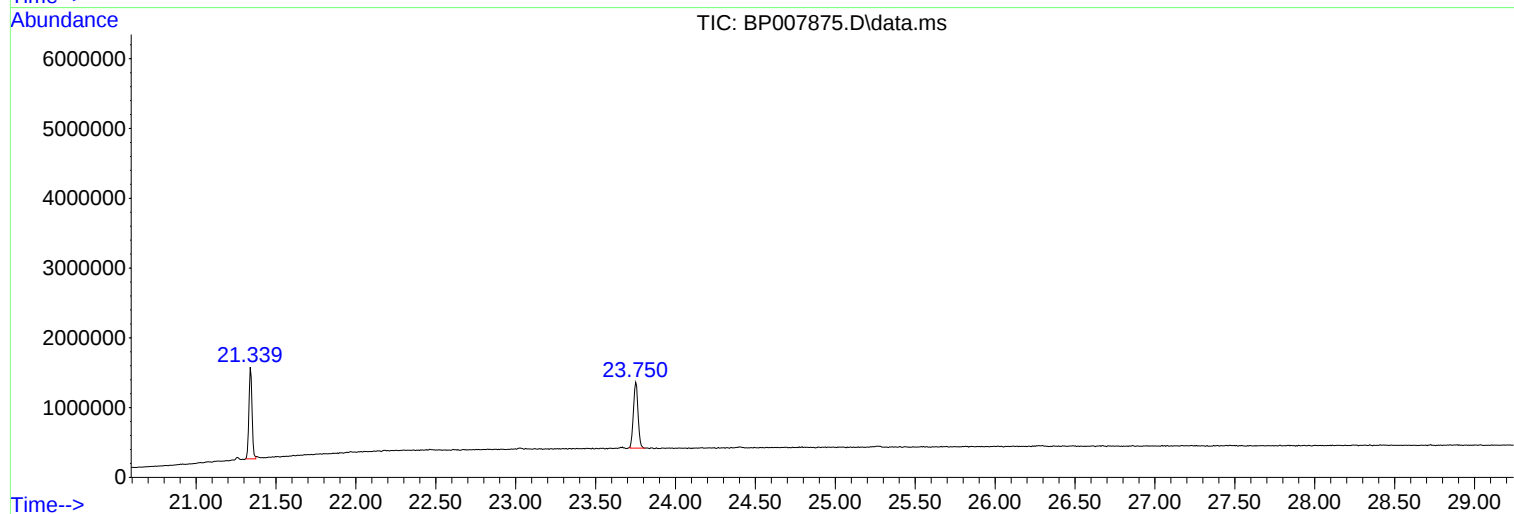
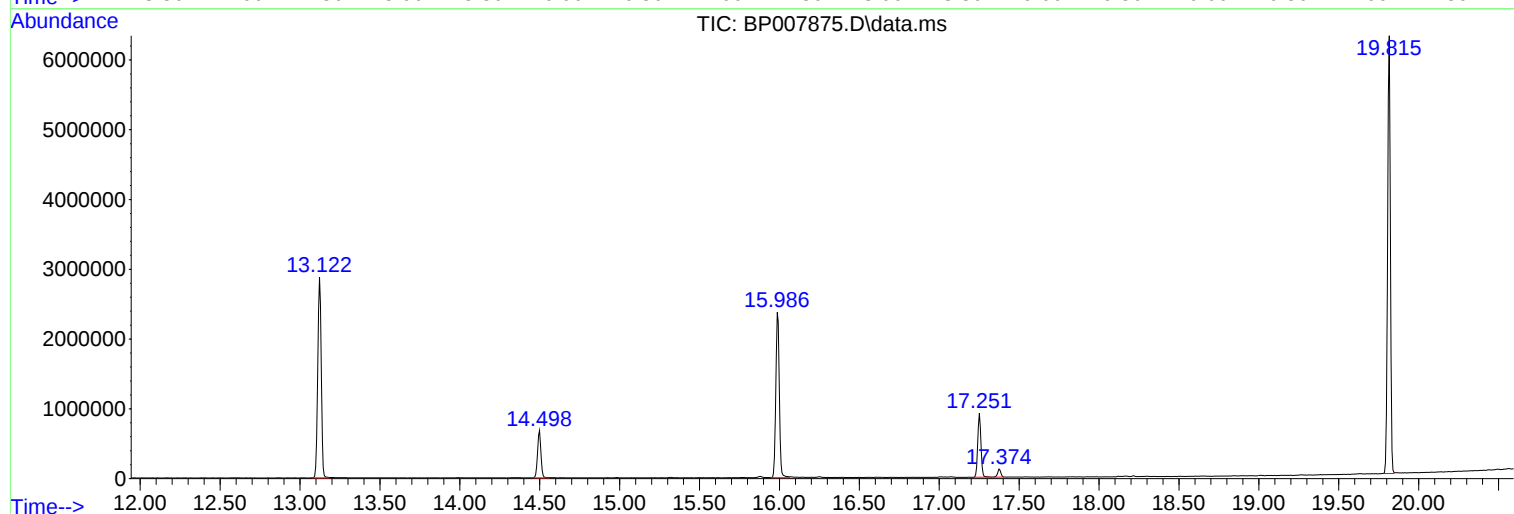
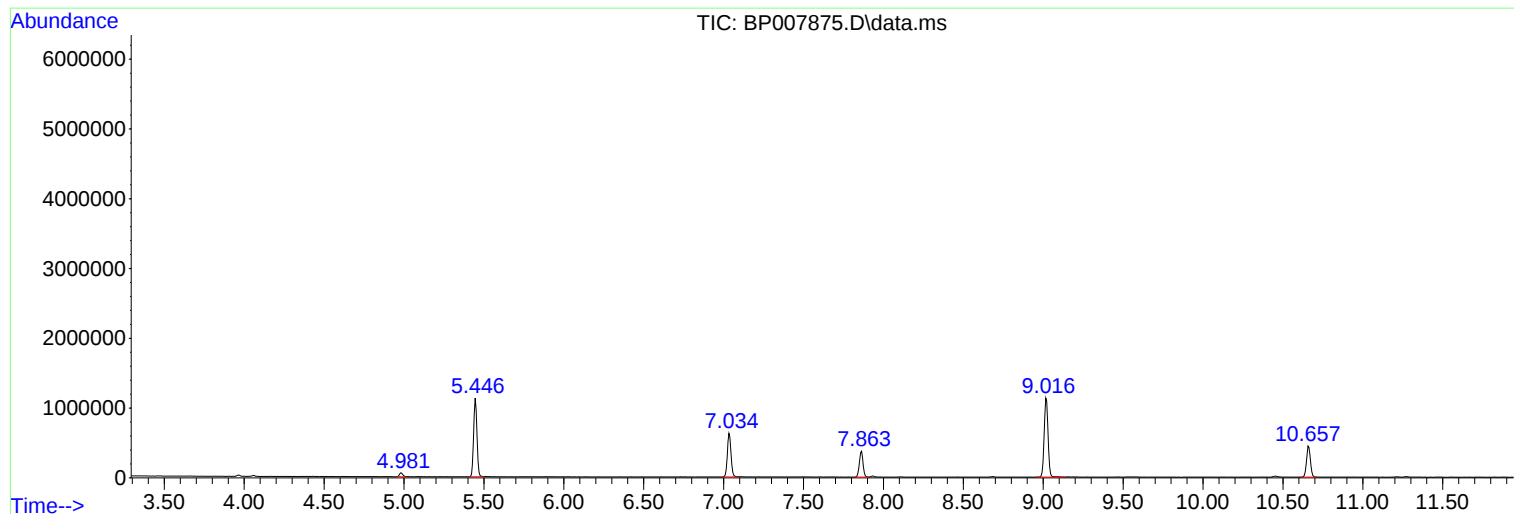
Sum of corrected areas: 27469581

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP110221.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P



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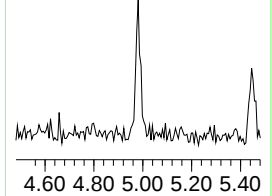
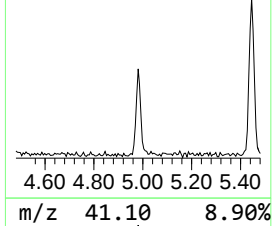
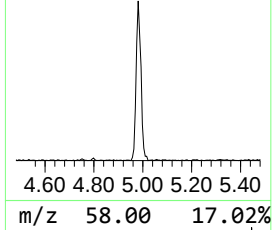
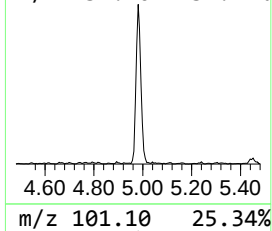
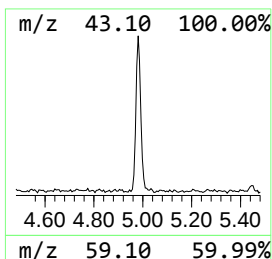
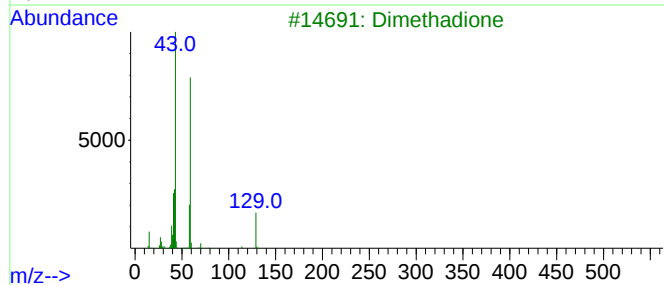
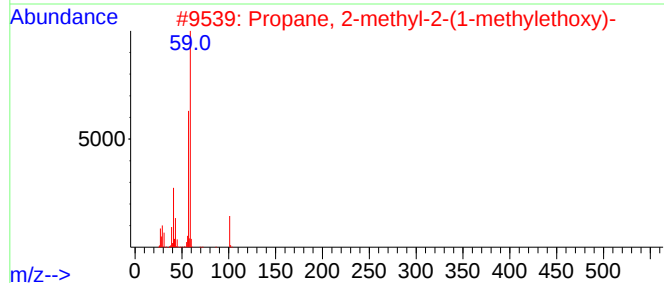
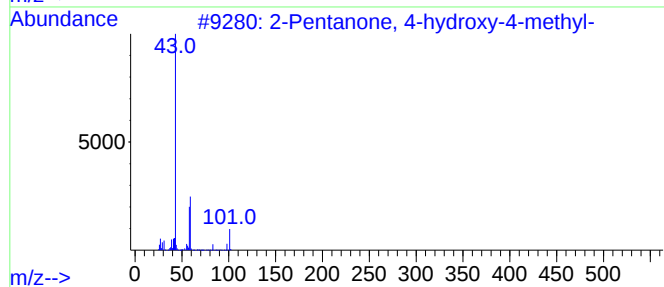
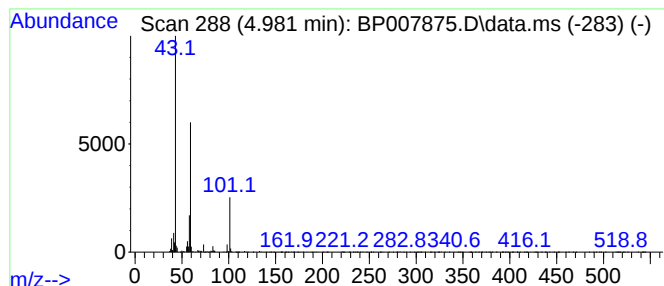
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TIC Library : C:\Database\NIST20.L
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.981	2.69 ng	79067	1,4-Dichlorobenzene-d4	7.863

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	59		
2	Propane, 2-methyl-2-(1-methyleth...	116	C7H16O	017348-59-3	50		
3	Dimethadione	129	C5H7NO3	000695-53-4	40		
4	Hydrazine, 1,1-bis(1-methylethyl)-	116	C6H16N2	000921-14-2	40		
5	(+)-4-Amino-4,5-dihydro-2(3H)-f...	101	C4H7NO2	016504-58-8	38		



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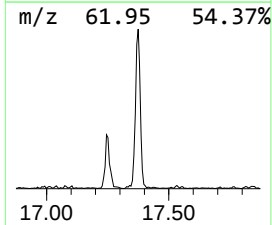
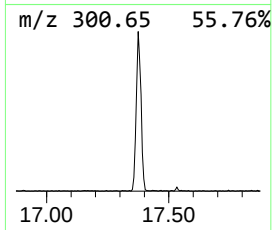
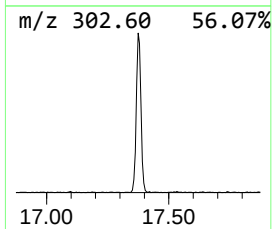
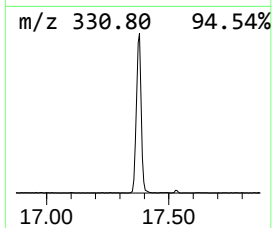
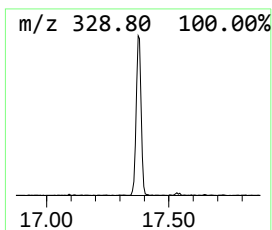
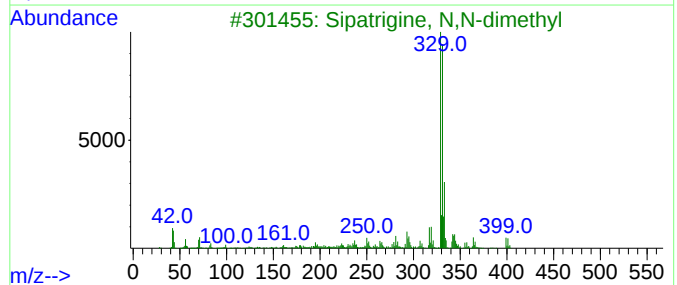
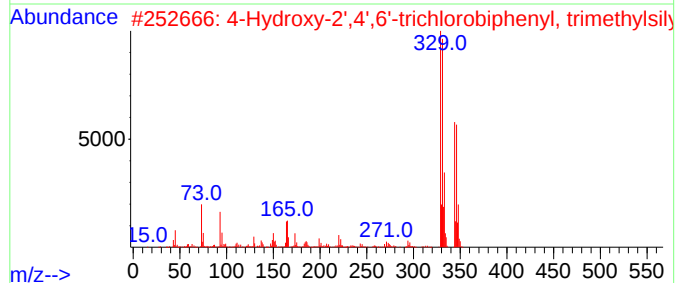
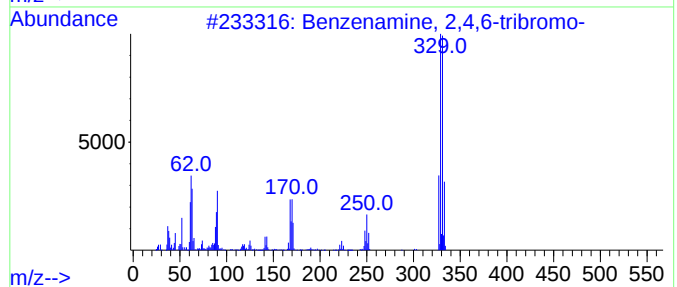
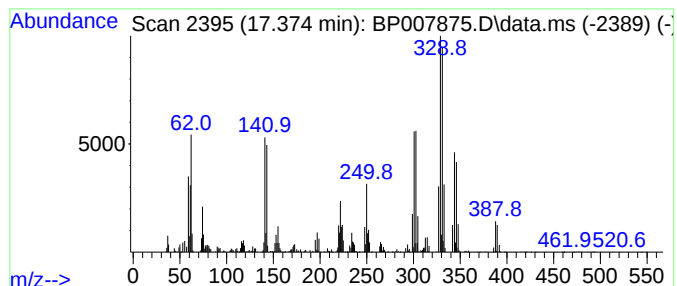
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Peak Number 2 Benzenamine, 2,4,6-tribromo- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD		R.T.	
17.374	2.54 ng	160804	Phenanthrene-d10		17.251	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Benzenamine, 2,4,6-tribromo-		327	C6H4Br3N	000147-82-0	64
2	4-Hydroxy-2',4',6'-trichlorobiph...		344	C15H15Cl3OSi	1000447-71-9	25
3	Sipatrigine, N,N-dimethyl		399	C17H20Cl3N5	1000498-90-8	22
4	Allyl 2,4,6-tribromophenyl ether		368	C9H7Br3O	003278-89-5	22
5	1,2,4-1H-Triazole, 1-(3,5-dichlo...		329	C12H13Cl2N5S	1000266-38-9	18



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-...	4.981	2.7	ng	79067	1	7.863	587526	20.0
Benzenamine, 2,...	17.374	2.5	ng	160804	4	17.251	1264440	20.0