

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051021\
 Data File : BP005471.D
 Acq On : 10 May 2021 16:36
 Operator : CG/JU
 Sample : M2295-01
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 DRUM-A-B-C

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-BP050521.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BP005471.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	4.834	290	297	304	rBV	332582	419201	48.27%	10.591%
2	5.293	370	375	384	rBV	79652	115981	13.35%	2.930%
3	6.893	640	647	662	rBV	139463	228386	26.30%	5.770%
4	7.705	778	785	793	rBV	72952	110572	12.73%	2.794%
5	8.869	976	983	995	rBV	207434	332545	38.29%	8.402%
6	10.493	1253	1259	1267	rBV	67198	115233	13.27%	2.911%
7	12.969	1672	1680	1692	rBV	493094	713586	82.17%	18.029%
8	13.822	1819	1825	1832	rBV	6323	10188	1.17%	0.257%
9	14.351	1908	1915	1922	rBV	93402	138645	15.96%	3.503%
10	15.845	2160	2169	2184	rBV	78068	136907	15.76%	3.459%
11	17.104	2376	2383	2393	rBV	101816	161954	18.65%	4.092%
12	19.669	2813	2819	2828	rBV	723436	868448	100.00%	21.942%
13	20.927	3030	3033	3038	rVB	14075	18967	2.18%	0.479%
14	20.980	3038	3042	3048	rBV	46665	62640	7.21%	1.583%
15	21.198	3074	3079	3083	rBV	155677	200691	23.11%	5.071%
16	22.916	3368	3371	3375	rBV6	13824	26901	3.10%	0.680%
17	23.510	3466	3472	3483	rVB2	144116	297165	34.22%	7.508%

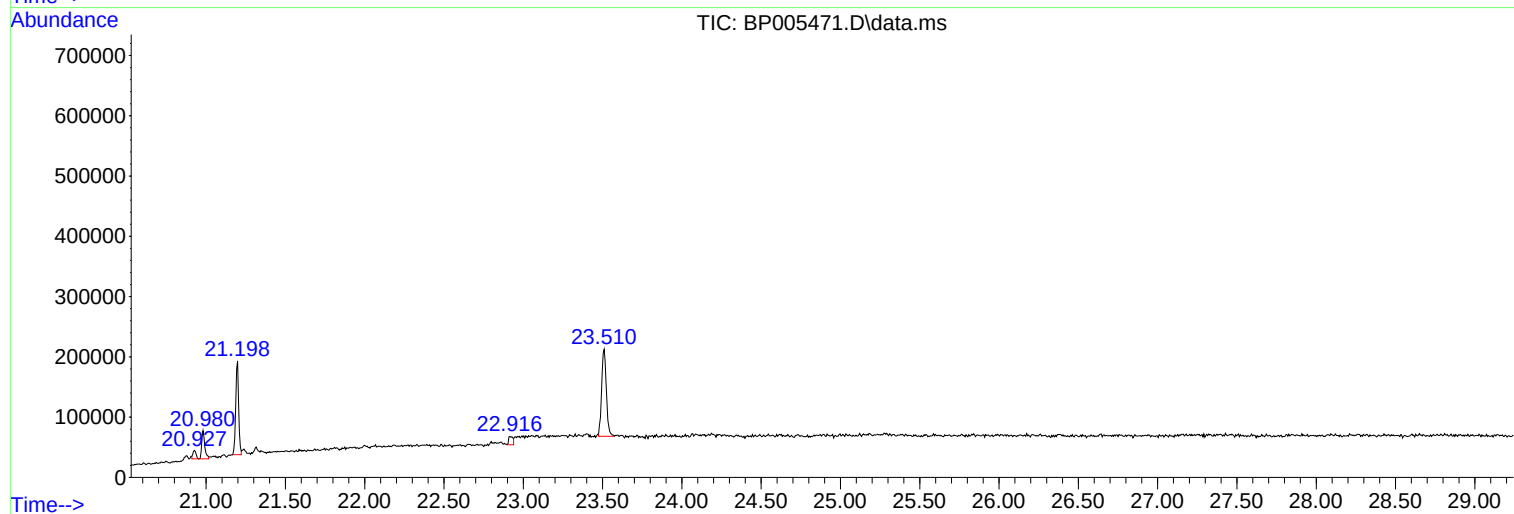
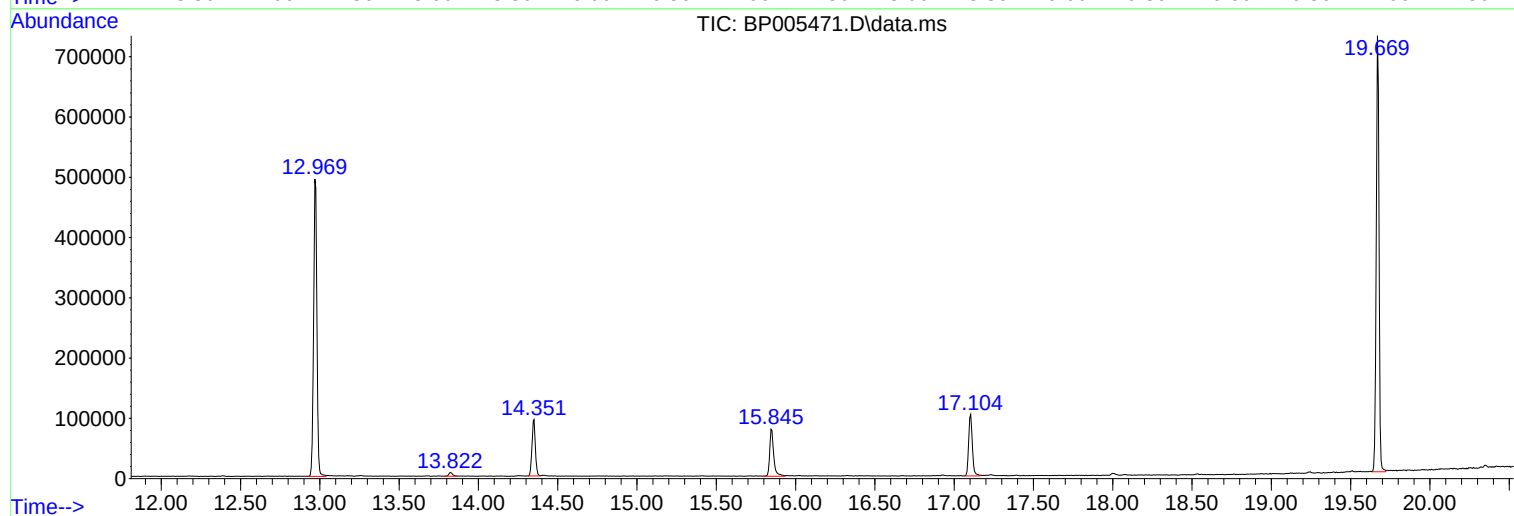
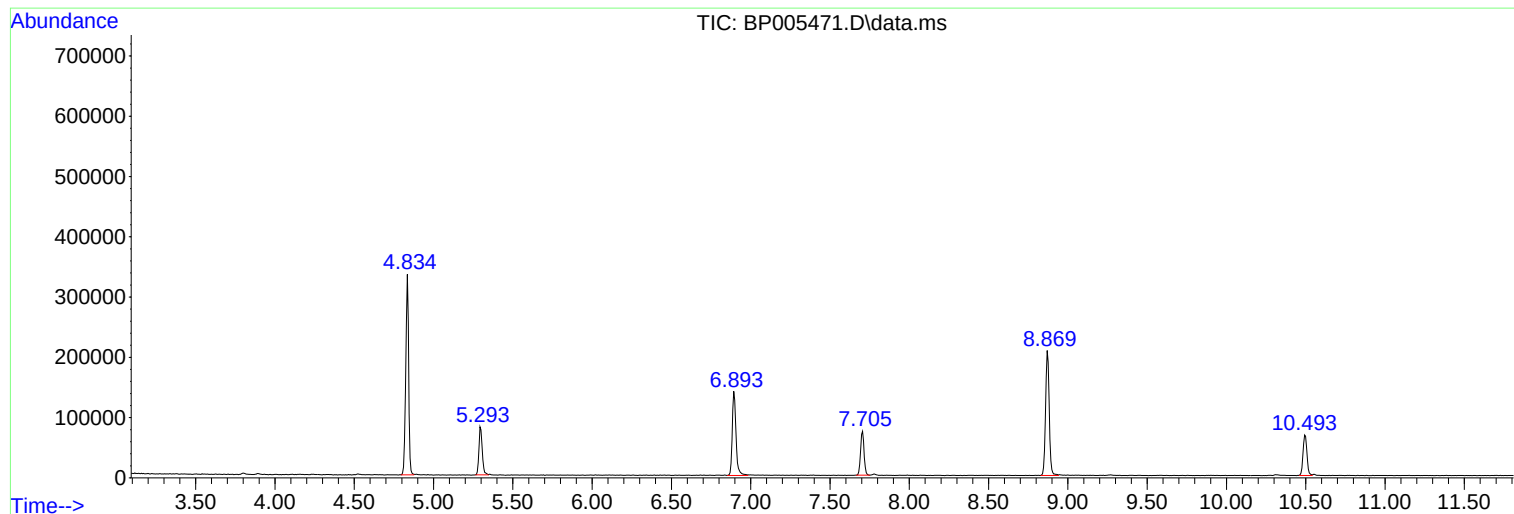
Sum of corrected areas: 3958010

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP050521.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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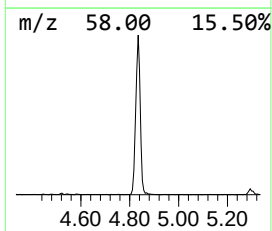
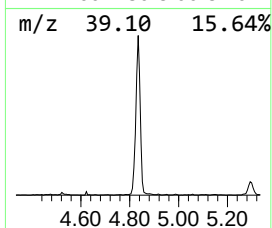
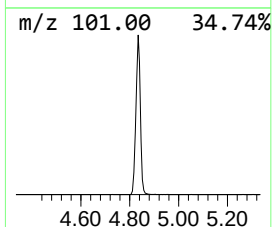
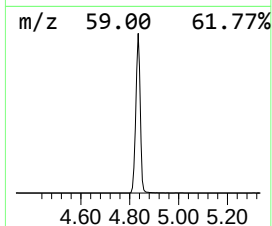
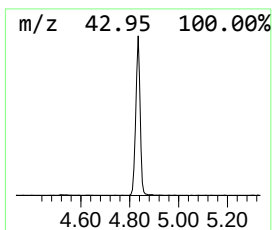
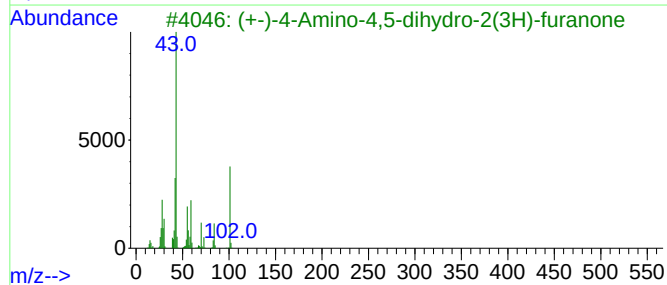
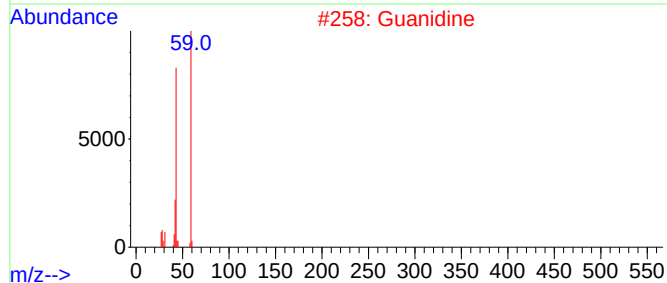
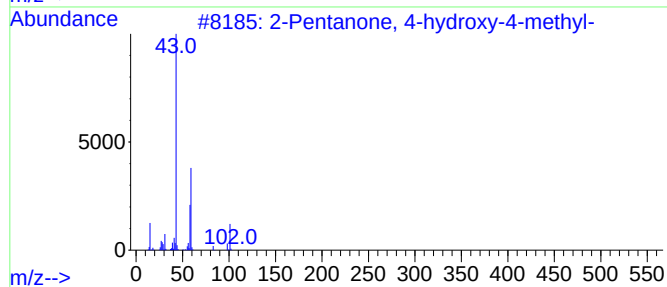
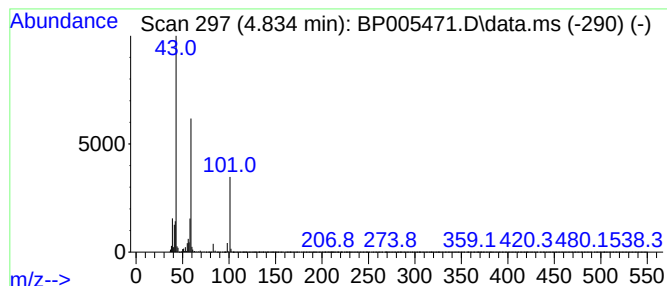
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP050521.M
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TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.834	75.82 ng	419201	1,4-Dichlorobenzene-d4	7.705

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	38		
2	Guanidine	59	CH5N3	000113-00-8	37		
3	(+)-4-Amino-4,5-dihydro-2(3H)-f...	101	C4H7NO2	016504-58-8	23		
4	2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	23		
5	Butane	58	C4H10	000106-97-8	22		



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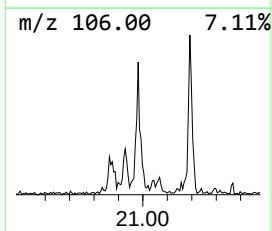
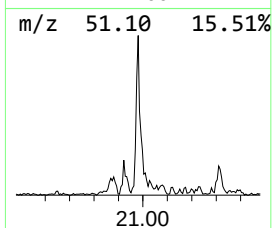
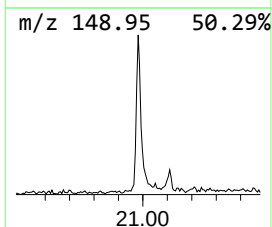
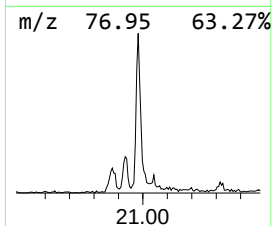
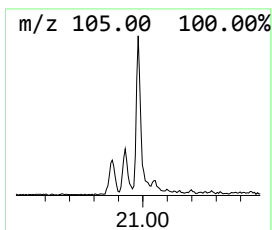
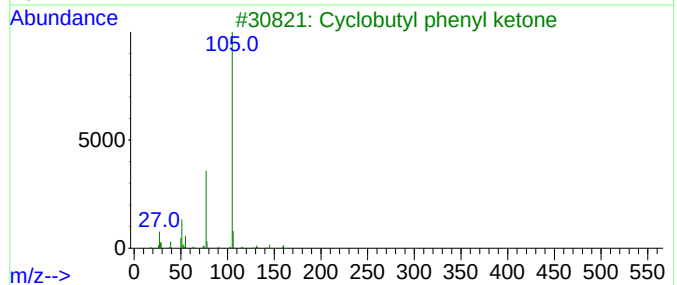
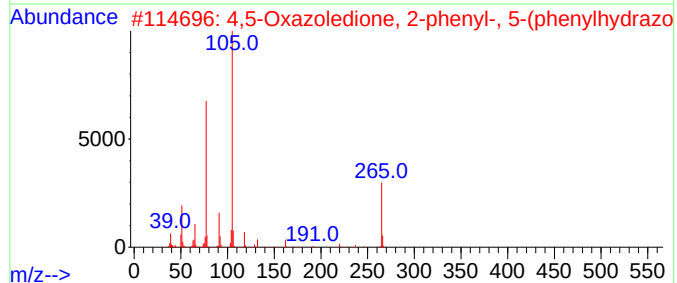
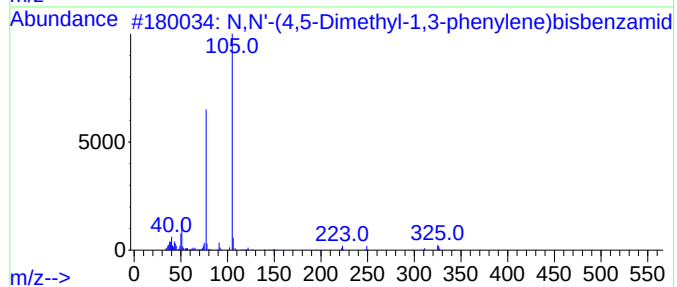
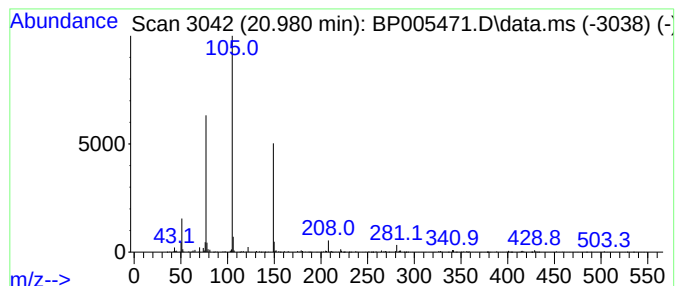
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TIC Integration Parameters: LSCINT.P

Peak Number 2 unknown20.98 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.980	6.24 ng	62640	Chrysene-d12	21.198

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			N,N'-(4,5-Dimethyl-1,3-phenylene...	344	C22H20N2O2	047507-95-9	38
2			4,5-Oxazolidione, 2-phenyl-, 5-(...	265	C15H11N3O2	1000337-82-6	38
3			Cyclobutyl phenyl ketone	160	C11H12O	005407-98-7	38
4			N-Methoxy-N-methylbenzamide	165	C9H11NO2	006919-61-5	38
5			Benzeneacetic acid, .alpha.-oxo-...	164	C9H8O3	015206-55-0	38



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-...	4.834	75.8	ng	419201	1	7.705	110572	20.0
unknown20.98	20.980	6.2	ng	62640	5	21.198	200691	20.0