

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051921\
Data File : BP005614.D
Acq On : 19 May 2021 16:36
Operator : CG/JU
Sample : PB136538BL
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
PB136538BL

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: BP005614.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	4.793	285	290	301	rBV2	5522	10236	1.04%	0.315%
2	5.252	362	368	386	rBV	171125	278818	28.29%	8.592%
3	6.857	634	641	658	rBV	136157	250979	25.46%	7.734%
4	7.657	769	777	788	rBV	35450	58616	5.95%	1.806%
5	8.828	969	976	992	rBV	90970	165198	16.76%	5.090%
6	10.451	1245	1252	1270	rBV2	31480	64489	6.54%	1.987%
7	12.928	1666	1673	1685	rBV	278081	416011	42.21%	12.819%
8	14.310	1902	1908	1922	rBV2	56062	89995	9.13%	2.773%
9	15.816	2157	2164	2194	rBV	262081	449094	45.56%	13.839%
10	17.069	2371	2377	2387	rBV	69104	118528	12.03%	3.652%
11	19.645	2809	2815	2824	rBV	811336	985624	100.00%	30.371%
12	21.174	3070	3075	3084	rBV	126801	180552	18.32%	5.564%
13	23.468	3459	3465	3474	rVB	85708	177092	17.97%	5.457%

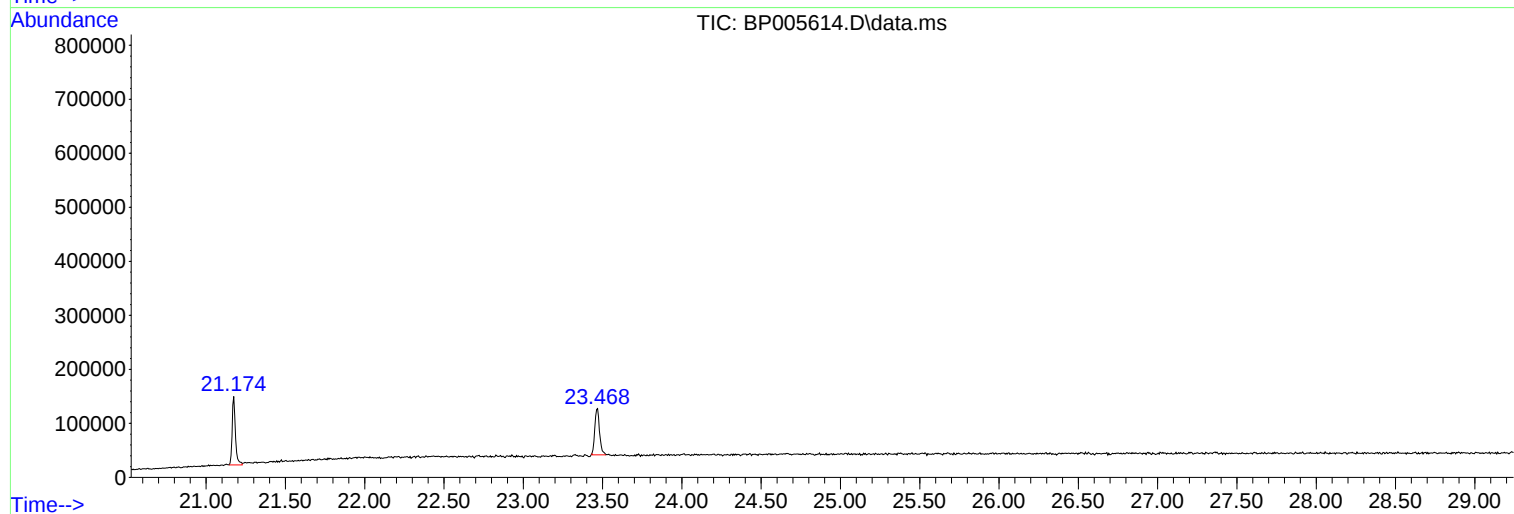
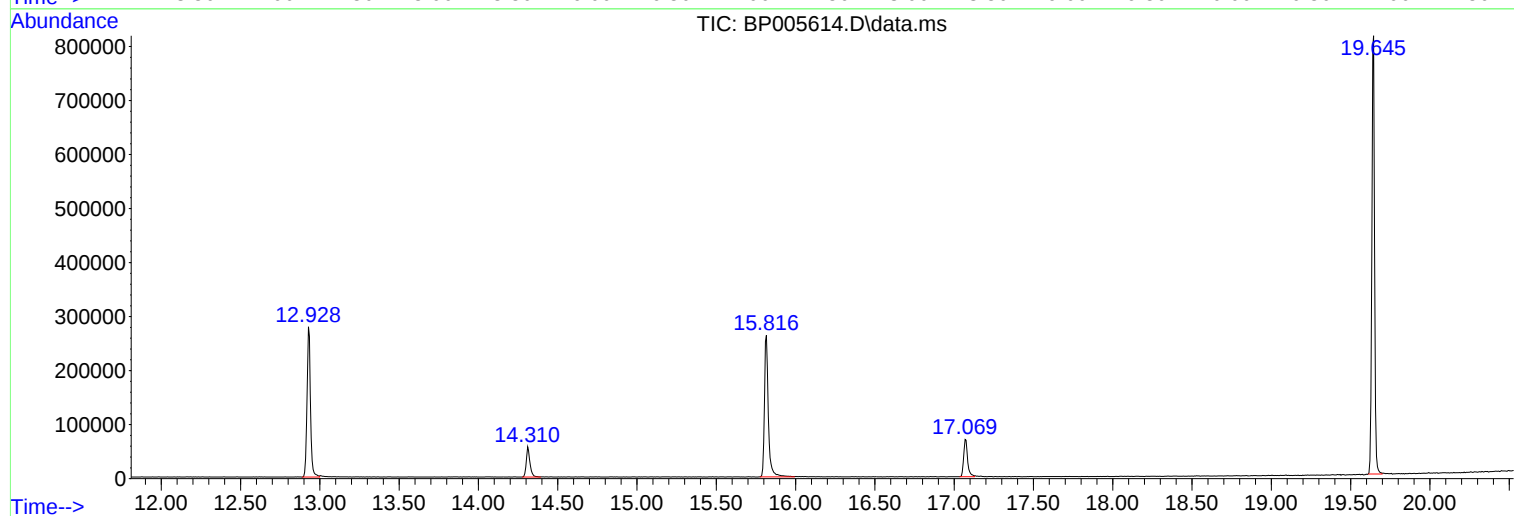
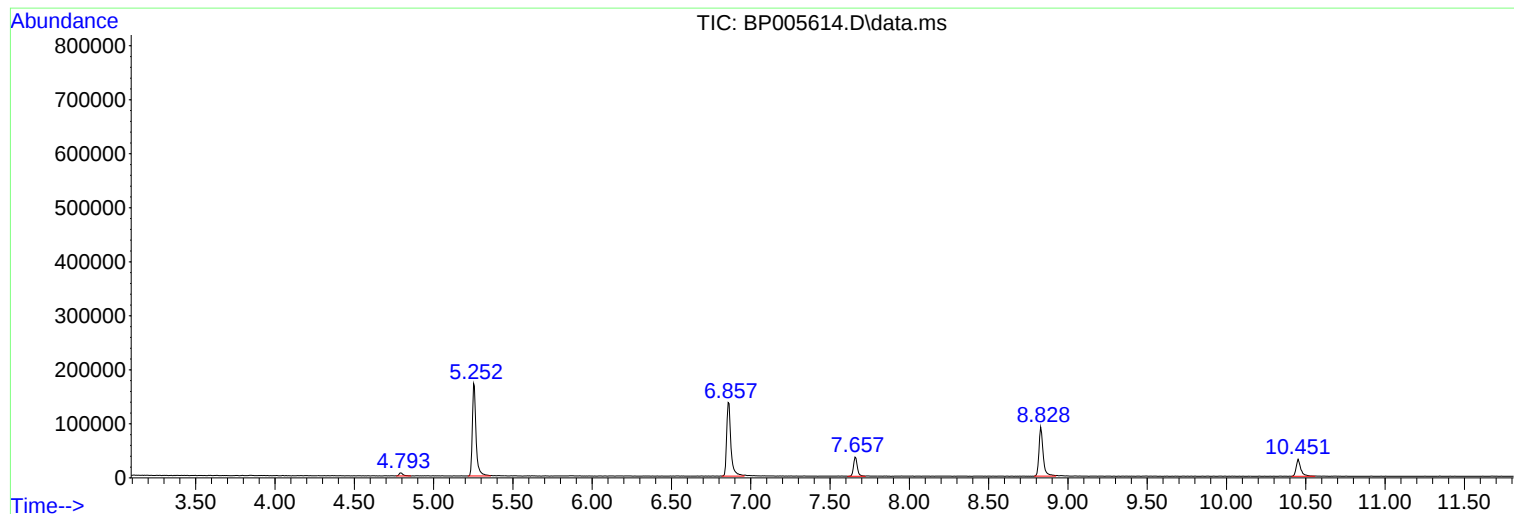
Sum of corrected areas: 3245232

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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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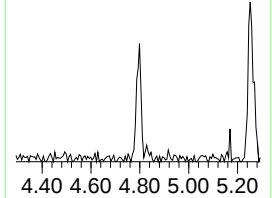
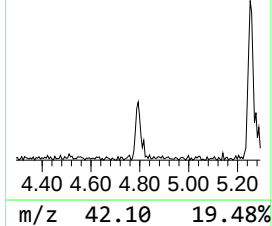
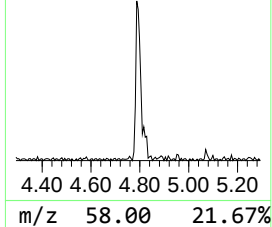
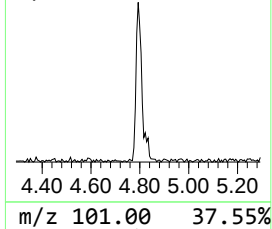
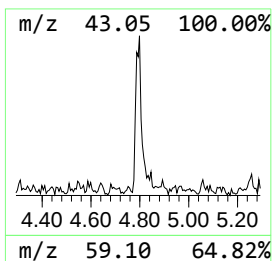
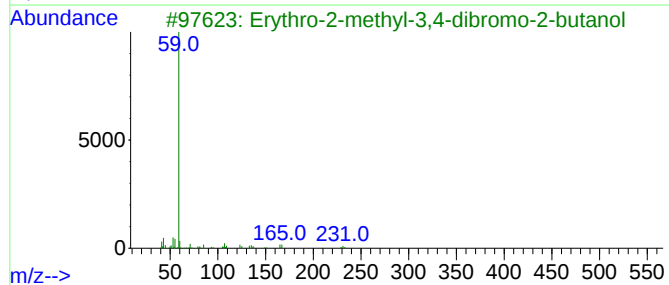
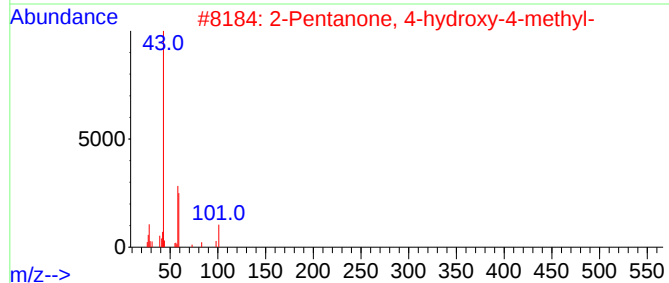
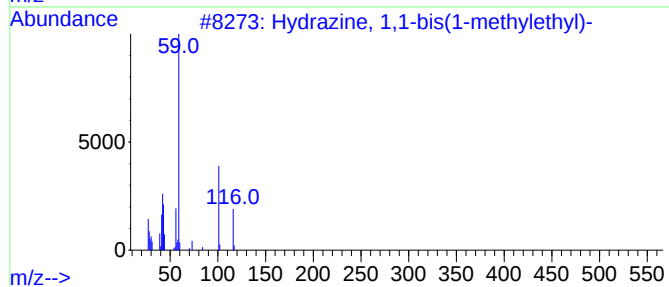
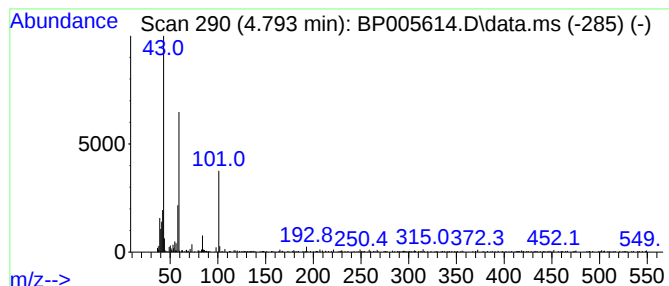
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Peak Number 1 unknown4.793 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.793	3.49 ng	10236	1,4-Dichlorobenzene-d4	7.663

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hydrazine, 1,1-bis(1-methylethyl)-	116	C6H16N2	000921-14-2	45
2			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	38
3			Erythro-2-methyl-3,4-dibromo-2-b...	244	C5H10Br2O	1000288-19-3	38
4			Nonanamide	157	C9H19NO	001120-07-6	38
5			2-Butanol, 3-methoxy-	104	C5H12O2	053778-72-6	32



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
unknown4.793	4.793	3.5	ng	10236	1	7.663	58616	20.0