

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP091922\
 Data File : BP011735.D
 Acq On : 19 Sep 2022 14:56
 Operator : CG/JU
 Sample : PB147234BL
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 PB147234BL

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-BP091422.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BP011735.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.034	342	348	357	rBV	285690	431207	2.14%	0.434%
2	5.499	419	427	444	rBV	7908208	11866685	58.86%	11.937%
3	7.099	691	699	718	rBV	7518873	11819445	58.62%	11.889%
4	7.940	834	842	853	rBV	1509967	2511462	12.46%	2.526%
5	9.105	1031	1040	1055	rBV	4379952	7177477	35.60%	7.220%
6	10.746	1309	1319	1334	rBV	1994225	3493507	17.33%	3.514%
7	13.204	1729	1737	1754	rBV	10190726	15235179	75.56%	15.325%
8	14.575	1963	1970	1985	rBV	2768891	4112880	20.40%	4.137%
9	16.063	2217	2223	2243	rBV	6282583	9225827	45.76%	9.280%
10	17.334	2432	2439	2451	rBV	3119311	4499963	22.32%	4.526%
11	19.880	2867	2872	2882	rBV	16185792	20162403	100.00%	20.281%
12	21.410	3127	3132	3138	rBV	3681183	4575569	22.69%	4.603%
13	23.863	3543	3549	3560	rVB2	2134238	4303212	21.34%	4.329%

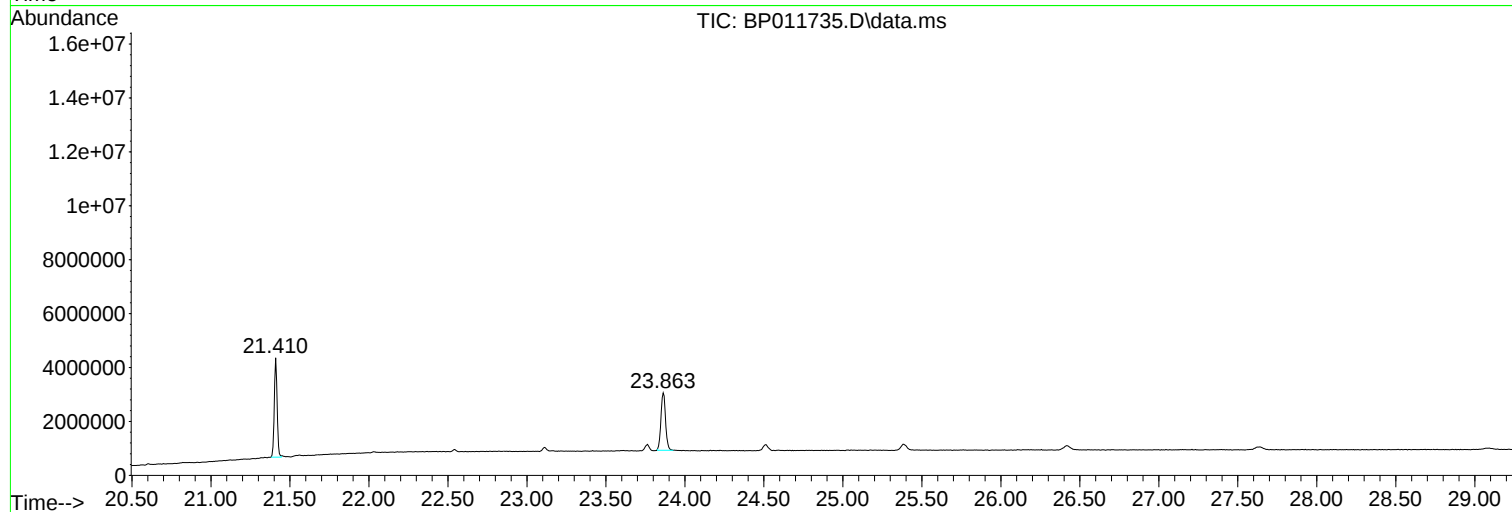
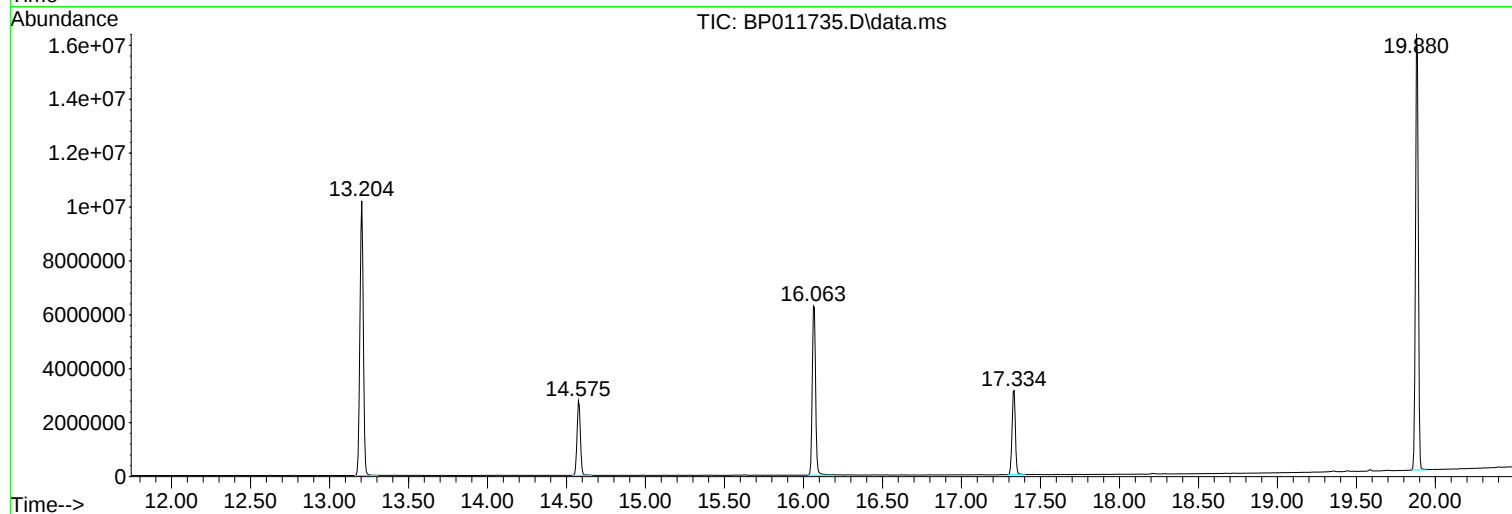
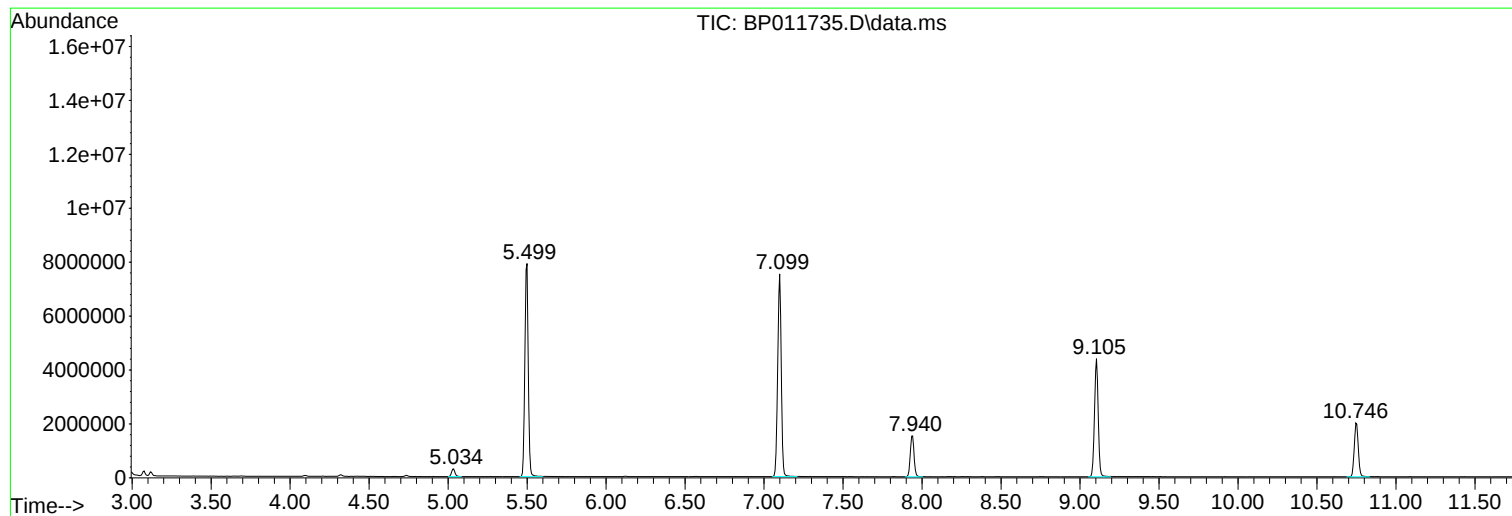
Sum of corrected areas: 99414816

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP091422.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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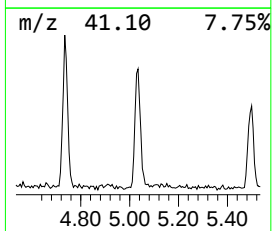
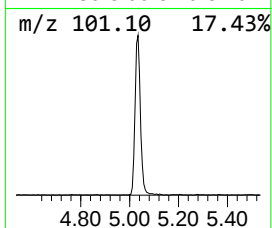
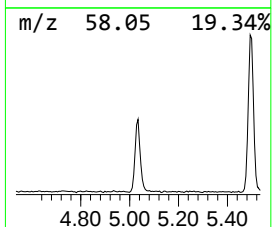
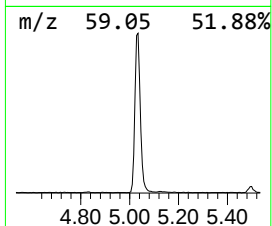
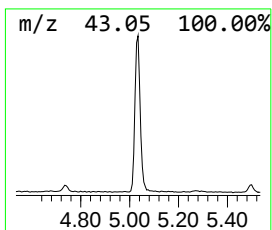
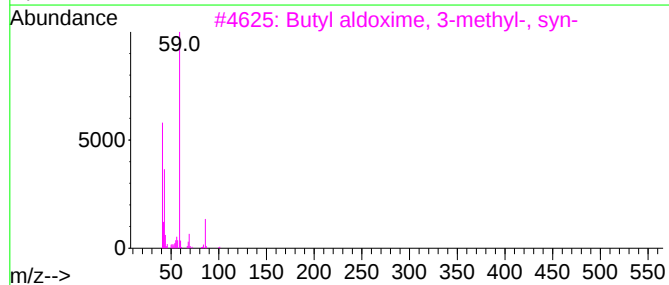
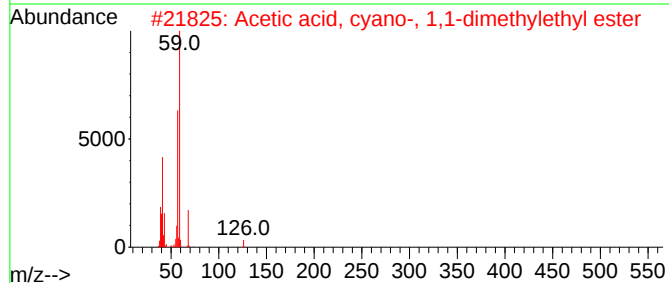
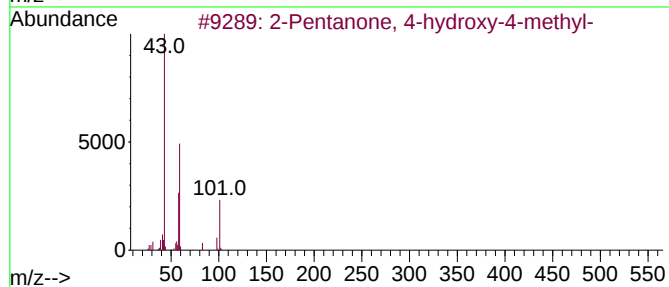
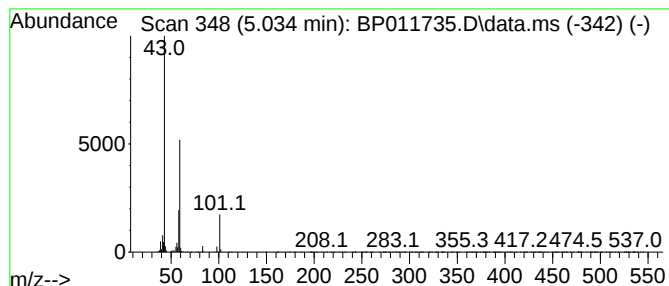
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TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD		R.T.	
5.034	3.43 ng	431207	1,4-Dichlorobenzene-d4		7.940	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-		116	C6H12O2	000123-42-2	83
2	Acetic acid, cyano-, 1,1-dimethy...		141	C7H11NO2	001116-98-9	23
3	Butyl aldoxime, 3-methyl-, syn-		101	C5H11NO	005780-40-5	17
4	2,3-Butanedione, monooxime		101	C4H7NO2	000057-71-6	12
5	Butane, 1-ethoxy-		102	C6H14O	000628-81-9	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-...	5.034	3.4	ng	431207	1	7.940	2511460	20.0