

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092121\
 Data File : BP007084.D
 Acq On : 21 Sep 2021 18:29
 Operator : CG/JU
 Sample : PB139176BL
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 PB139176BL

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-BP092121.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BP007084.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.011	322	327	335	rBV	377889	529520	3.46%	0.833%
2	5.475	399	406	422	rBV	4210978	6220881	40.65%	9.781%
3	7.069	669	677	693	rBV	3863737	6186140	40.42%	9.726%
4	7.905	811	819	828	rBV	765838	1242088	8.12%	1.953%
5	9.063	1008	1016	1034	rBV	2324079	3881496	25.36%	6.103%
6	10.705	1286	1295	1308	rBV	982764	1718583	11.23%	2.702%
7	13.157	1704	1712	1728	rBV	6560090	9551979	62.42%	15.019%
8	14.534	1937	1946	1958	rBV	1528491	2325492	15.20%	3.656%
9	16.016	2191	2198	2218	rBV	4935716	7148177	46.71%	11.239%
10	17.281	2406	2413	2426	rBV	1936730	2816884	18.41%	4.429%
11	19.839	2842	2848	2858	rBV	12204851	15302885	100.00%	24.061%
12	21.145	3066	3070	3078	rBV	120651	200498	1.31%	0.315%
13	21.363	3102	3107	3115	rBV	2477386	3214511	21.01%	5.054%
14	23.774	3511	3517	3526	rVB2	1562643	3261867	21.32%	5.129%

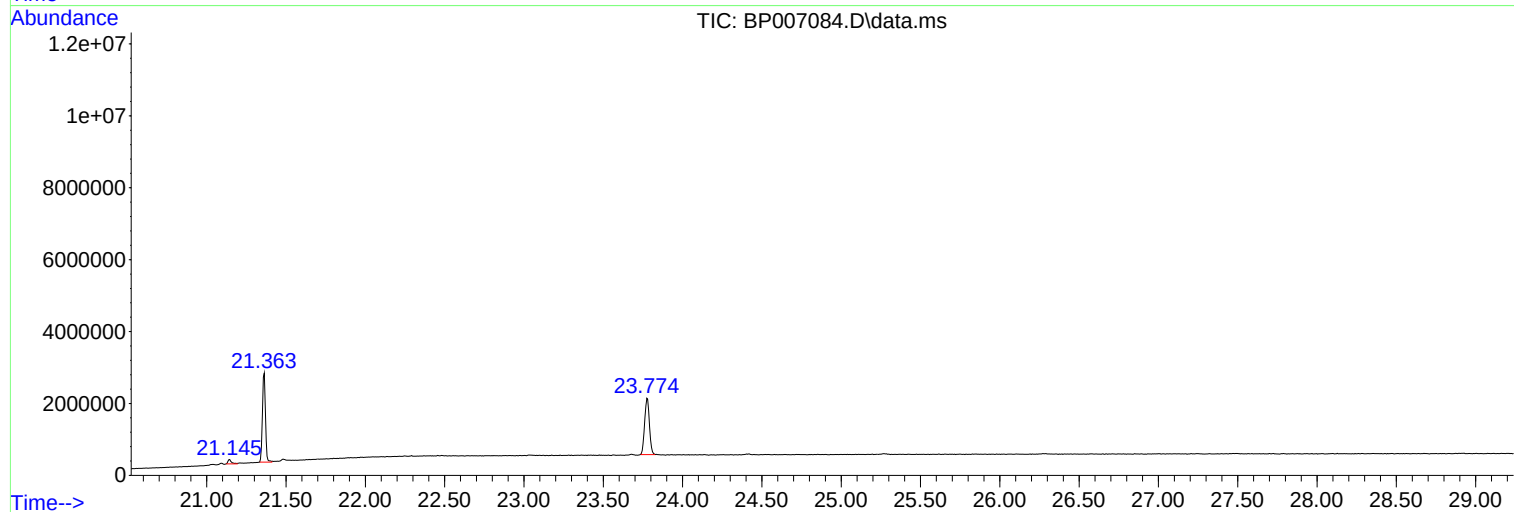
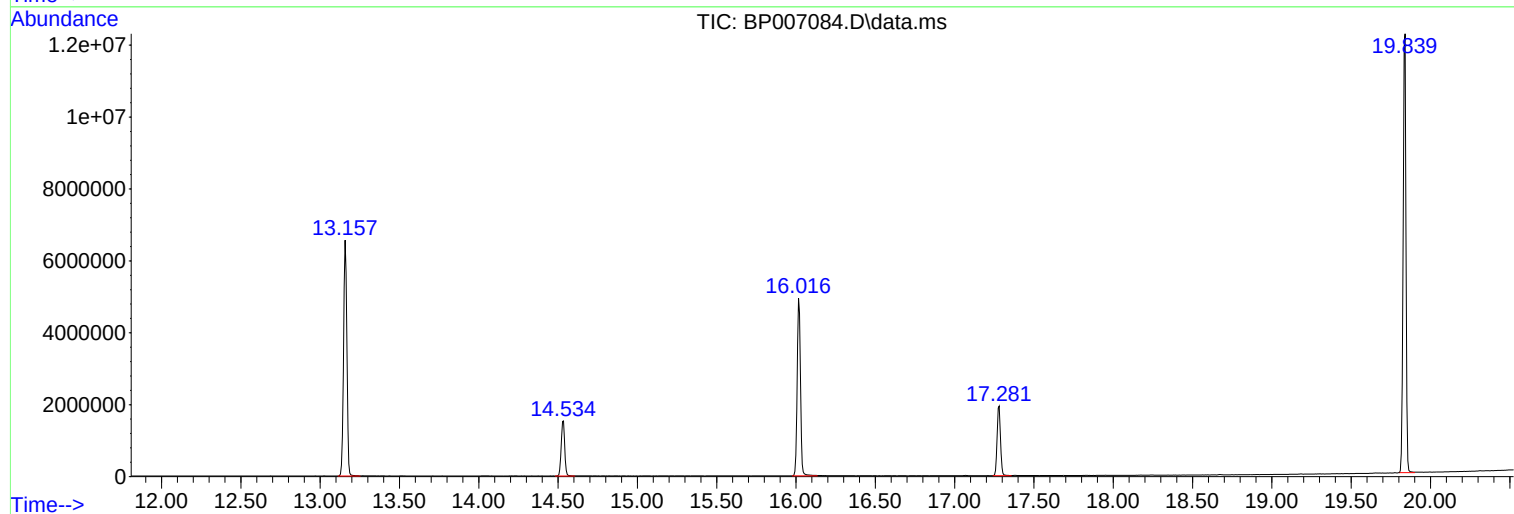
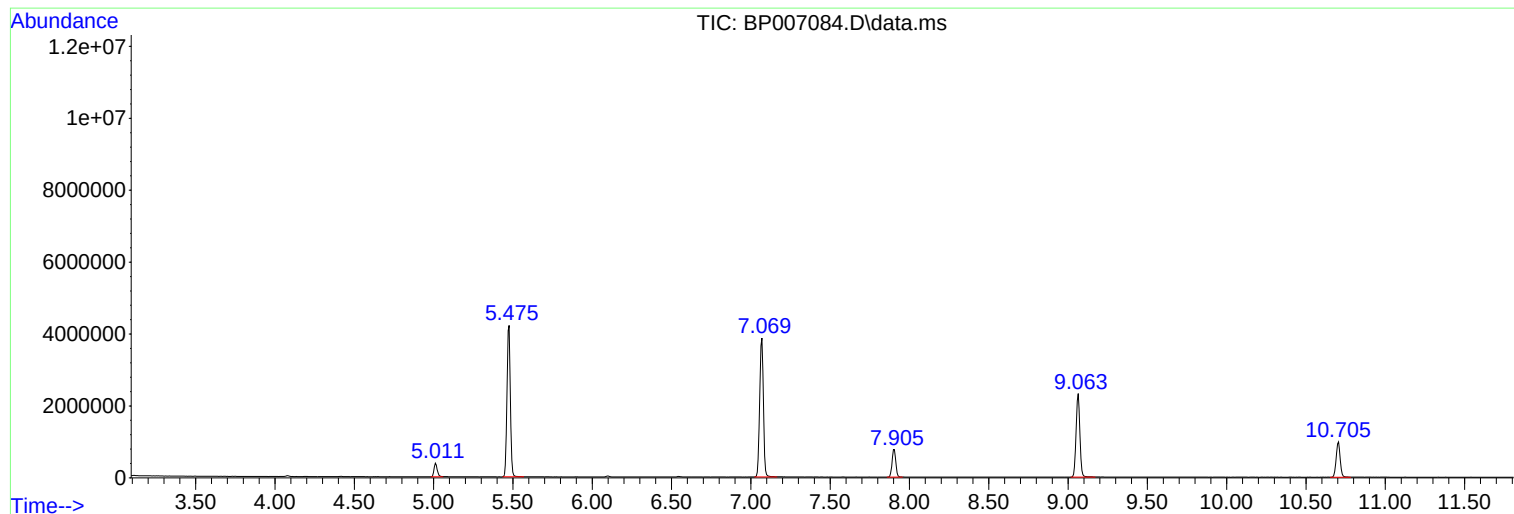
Sum of corrected areas: 63601001

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP092121.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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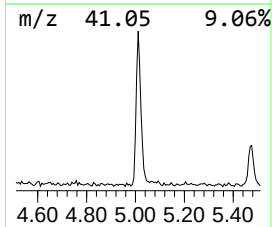
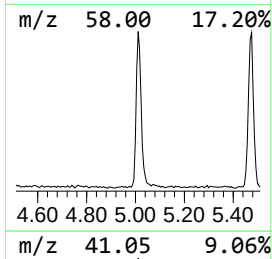
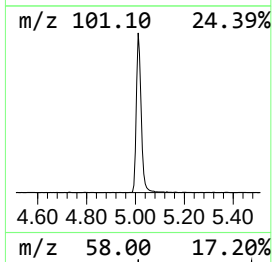
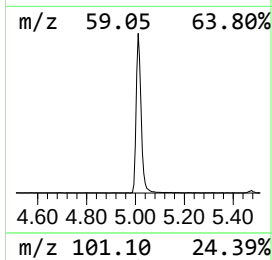
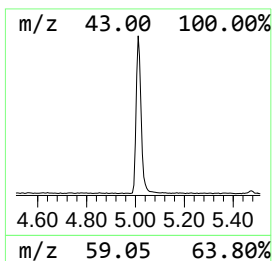
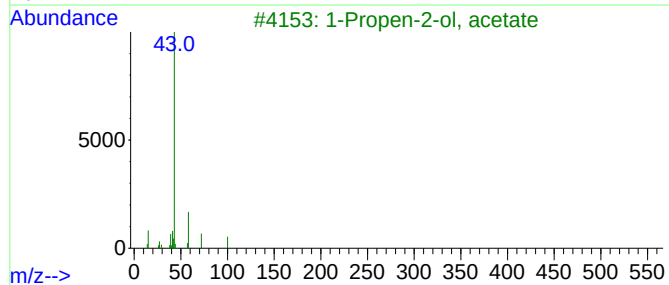
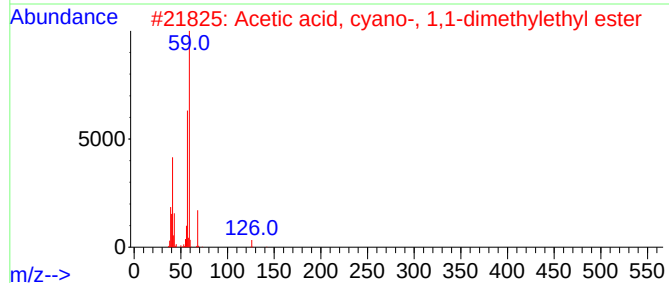
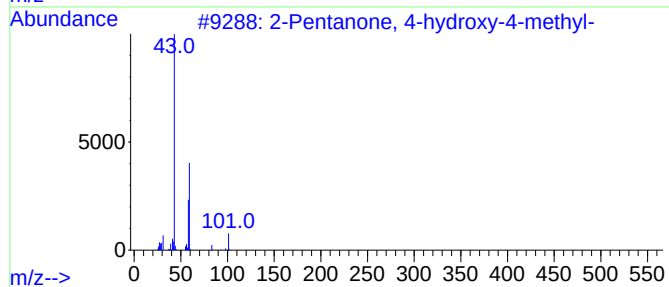
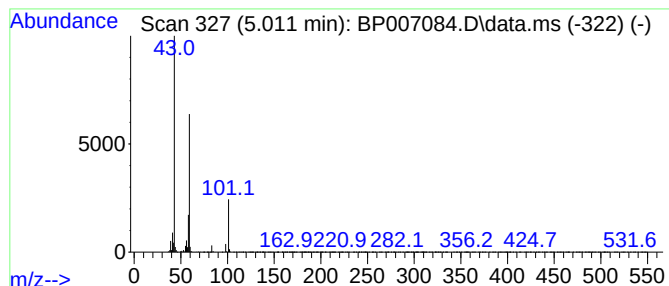
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.011	8.53 ng	529520	1,4-Dichlorobenzene-d4	7.905

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56		
2	Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	25		
3	1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	10		
4	Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9		
5	(+)-4-Amino-4,5-dihydro-2(3H)-f...	101	C4H7NO2	016504-58-8	9		



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-...	5.011	8.5	ng	529520	1	7.905	1242090	20.0