

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP091721\
 Data File : BP007025.D
 Acq On : 17 Sep 2021 14:24
 Operator : CG/JU
 Sample : PB139165BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 PB139165BL

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-BP091621.M
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BP007025.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.016	324	328	340	rBV	201890	296824	1.80%	0.421%
2	5.481	400	407	420	rBV	4800564	7090979	42.95%	10.048%
3	7.069	668	677	691	rBV	4404057	7003985	42.42%	9.924%
4	7.910	812	820	828	rBV	980971	1553207	9.41%	2.201%
5	9.069	1007	1017	1033	rBV	2615559	4335662	26.26%	6.143%
6	10.710	1287	1296	1310	rBV	1242015	2129893	12.90%	3.018%
7	13.163	1706	1713	1723	rBV	7084187	10444704	63.26%	14.800%
8	14.534	1937	1946	1960	rBV	1877261	2762671	16.73%	3.915%
9	16.022	2191	2199	2214	rBV	5150128	7512043	45.50%	10.644%
10	17.281	2406	2413	2425	rBV	2332648	3319360	20.10%	4.703%
11	19.839	2842	2848	2858	rBV	13282312	16510369	100.00%	23.394%
12	21.357	3101	3106	3116	rBV	2860131	3793973	22.98%	5.376%
13	23.774	3510	3517	3528	rVB2	1925568	3820534	23.14%	5.413%

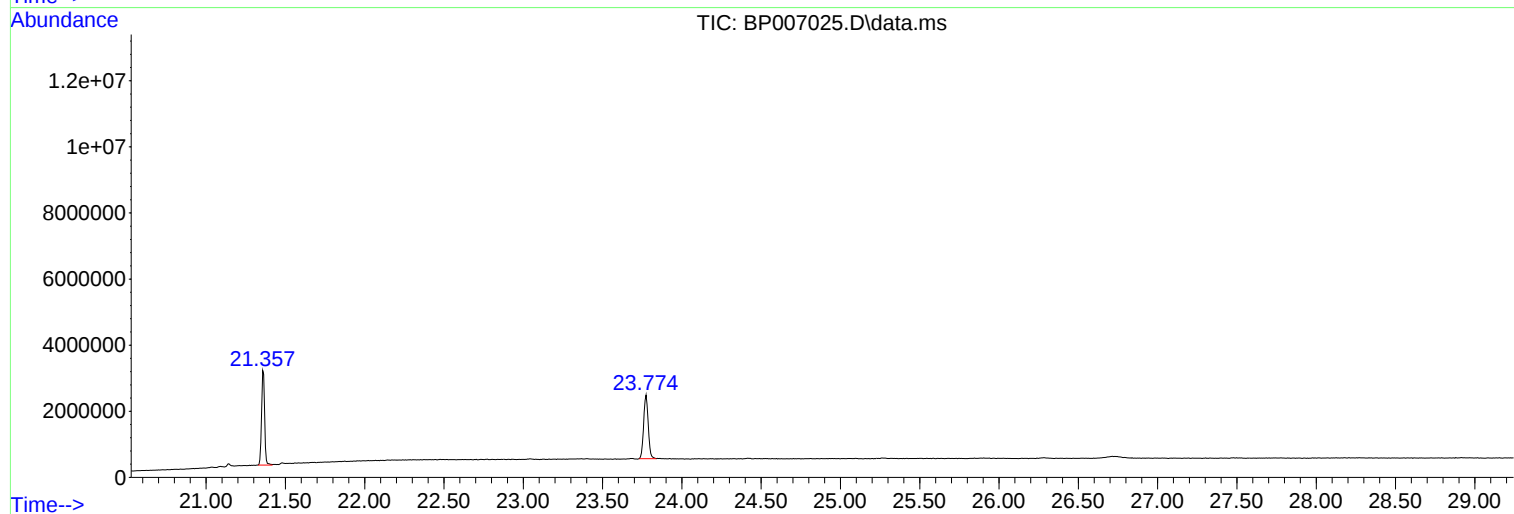
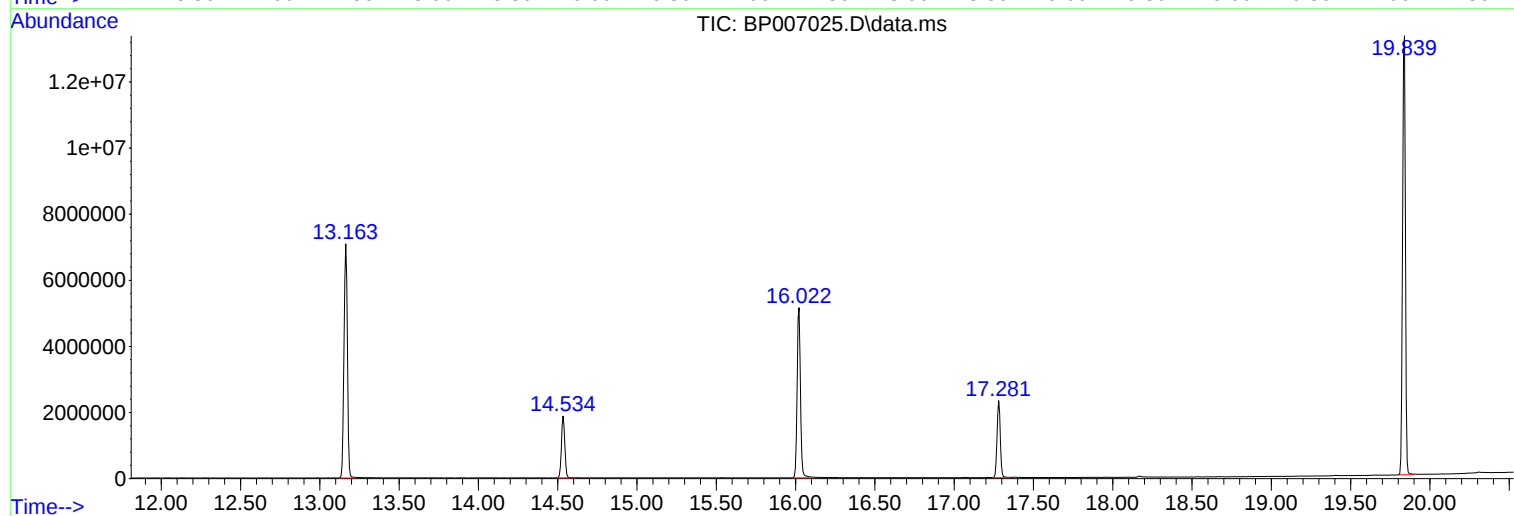
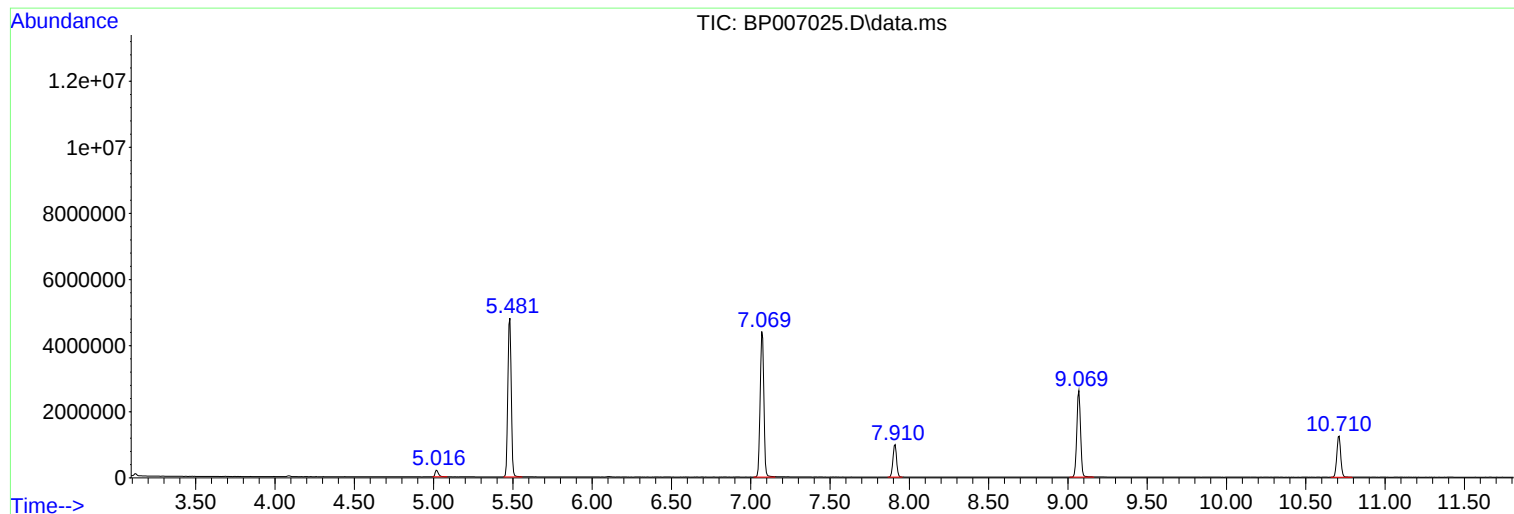
Sum of corrected areas: 70574204

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP091621.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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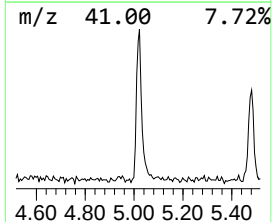
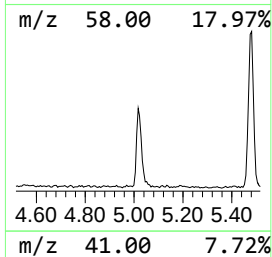
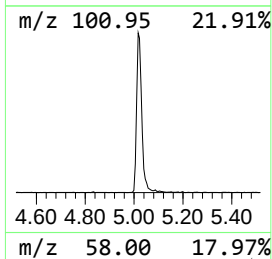
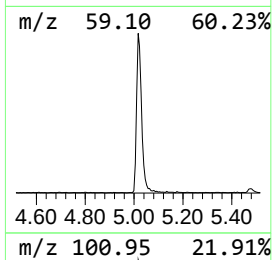
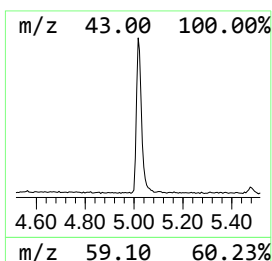
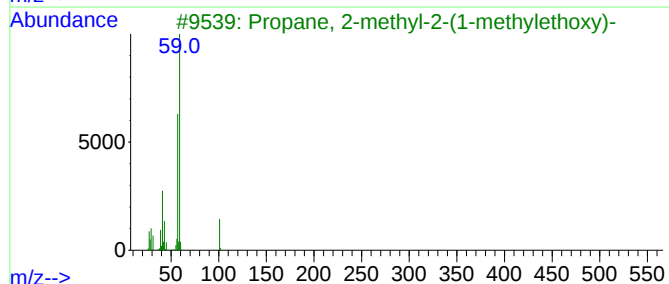
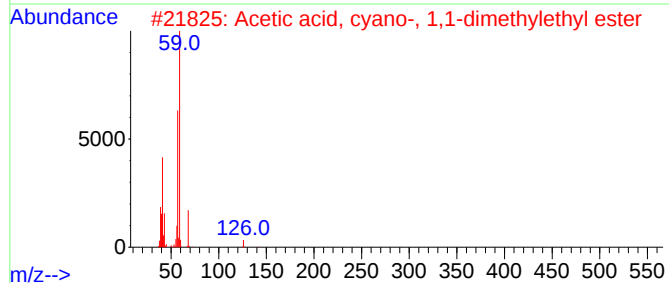
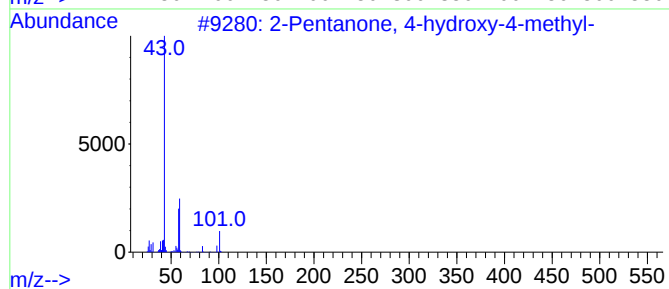
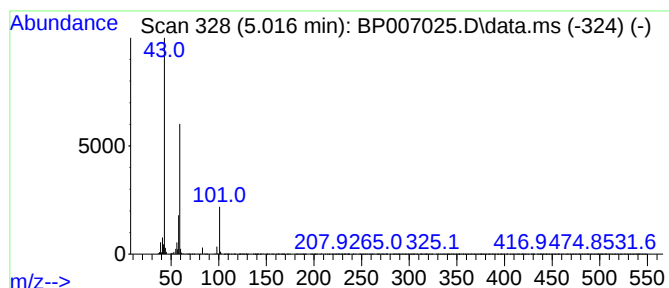
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD			R.T.
5.016	3.82 ng	296824	1,4-Dichlorobenzene-d4			7.910
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-		116	C6H12O2	000123-42-2	59
2	Acetic acid, cyano-, 1,1-dimethy...		141	C7H11NO2	001116-98-9	25
3	Propane, 2-methyl-2-(1-methyleth...		116	C7H16O	017348-59-3	23
4	2-Hexanol, 2-methyl-		116	C7H16O	000625-23-0	23
5	1-Propen-2-ol, acetate		100	C5H8O2	000108-22-5	10



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-...	5.016	3.8	ng	296824	1	7.910	1553210	20.0