

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP091622\
 Data File : BP011724.D
 Acq On : 16 Sep 2022 10:19
 Operator : CG/JU
 Sample : PB147206BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 PB147206BL

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: BP011724.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.075	10	15	19	rBV	296331	357705	1.63%	0.337%
2	5.028	341	347	359	rBV	558973	826631	3.78%	0.778%
3	5.493	419	426	441	rBV	8346898	12202174	55.73%	11.487%
4	7.093	690	698	725	rBV	7607886	12318734	56.26%	11.596%
5	7.934	834	841	851	rBV	1558566	2442035	11.15%	2.299%
6	9.105	1030	1040	1060	rBV	4469051	7570329	34.57%	7.126%
7	10.746	1312	1319	1334	rBV	2014938	3435106	15.69%	3.234%
8	13.204	1729	1737	1750	rBV	11010992	16699683	76.26%	15.721%
9	14.575	1963	1970	1980	rBV	2889353	4203127	19.20%	3.957%
10	16.063	2216	2223	2247	rBV	7477399	10576898	48.30%	9.957%
11	17.328	2432	2438	2452	rBV	3309171	4738357	21.64%	4.461%
12	19.880	2867	2872	2881	rBV	17407886	21896926	100.00%	20.613%
13	21.410	3127	3132	3138	rBV	3737198	4660121	21.28%	4.387%
14	23.863	3543	3549	3561	rVB2	2129751	4300568	19.64%	4.048%

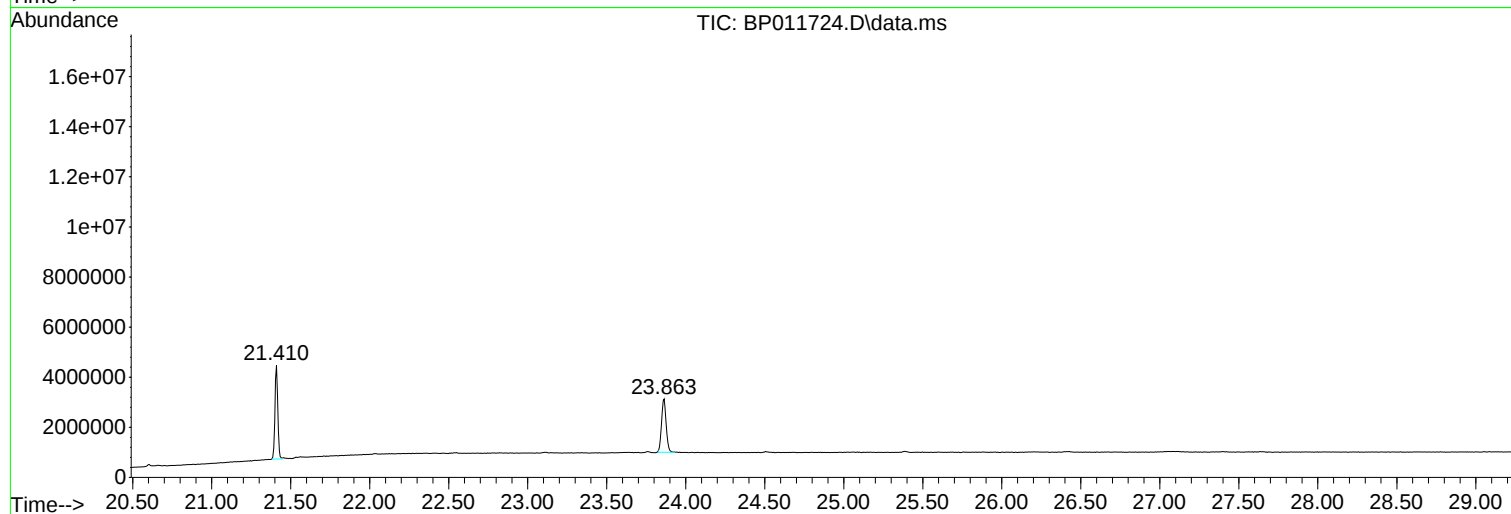
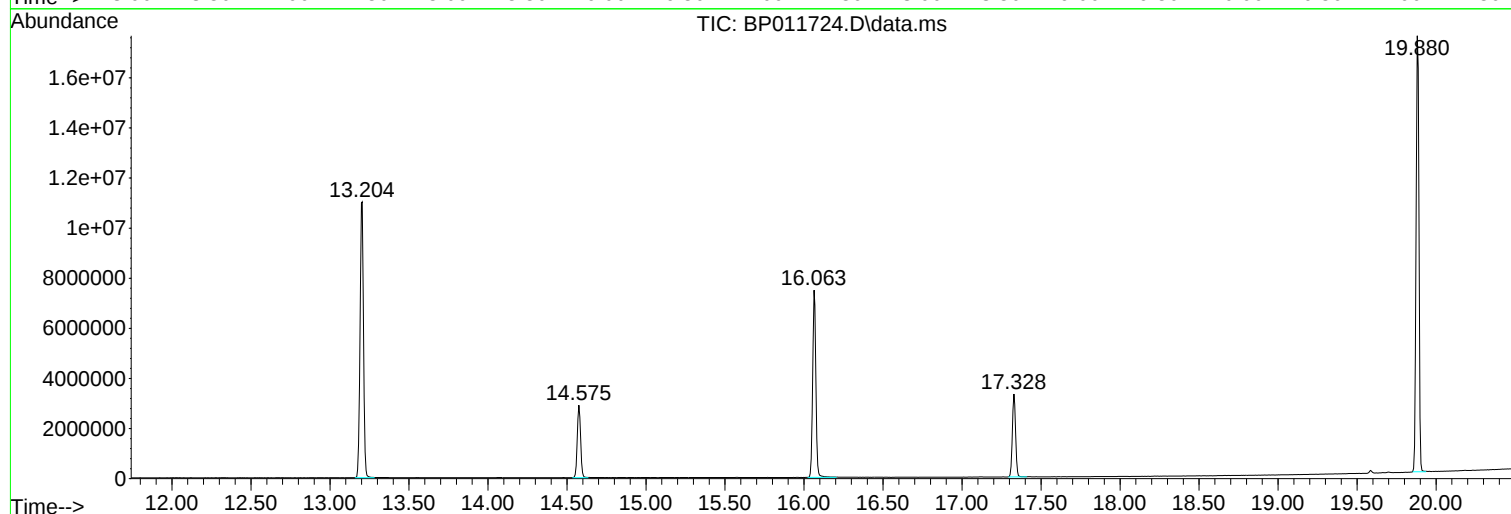
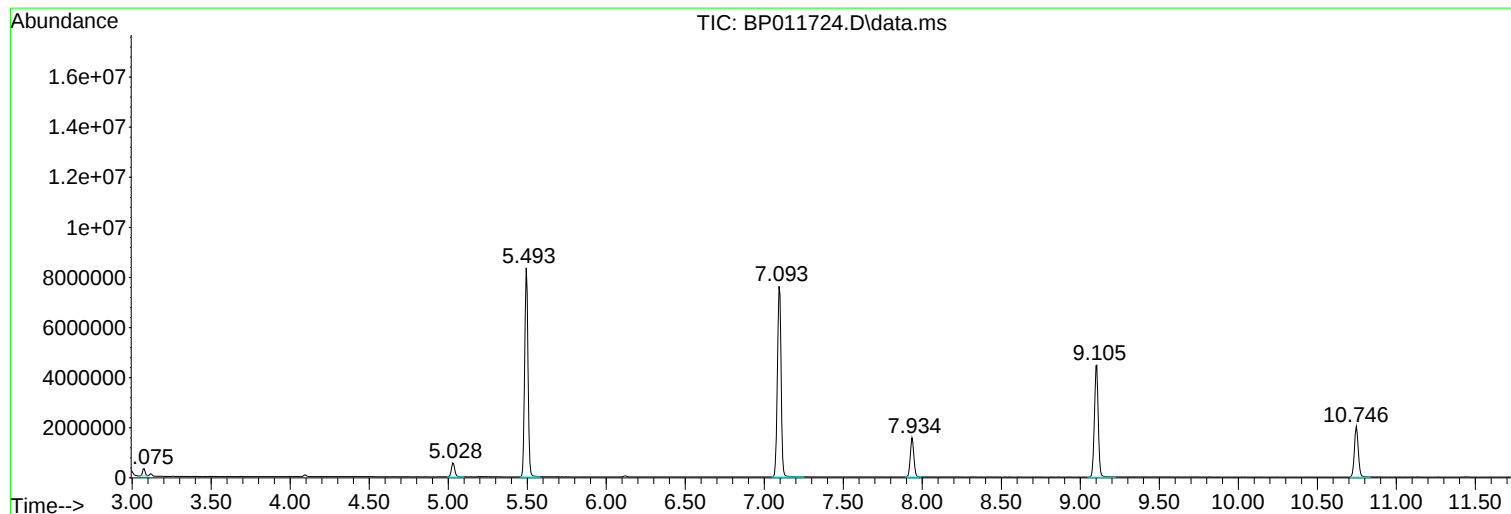
Sum of corrected areas: 106228394

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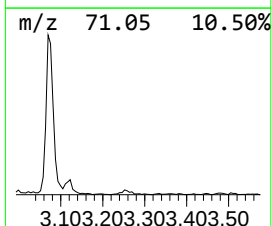
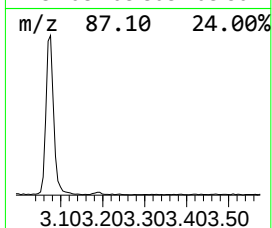
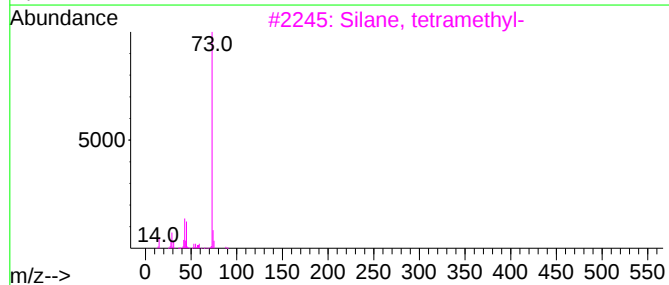
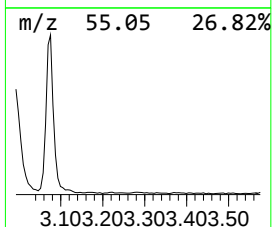
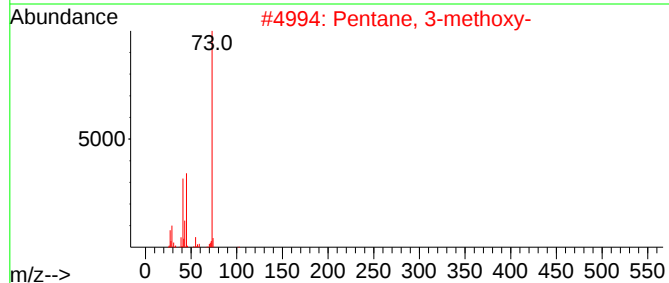
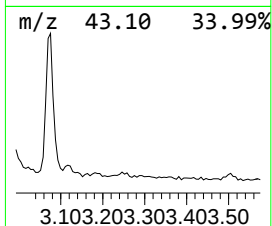
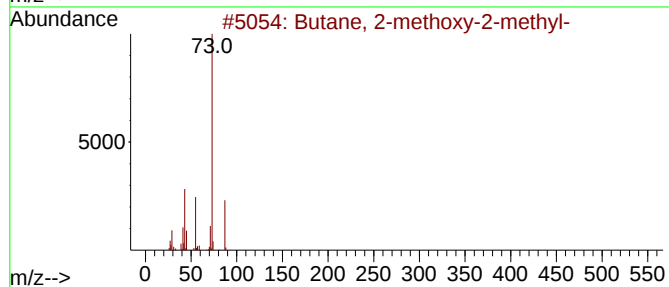
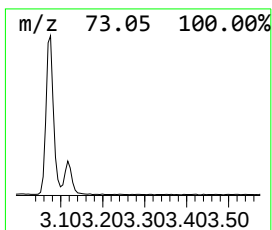
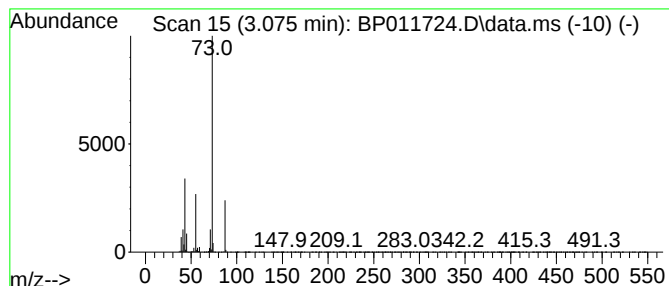
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP091422.M
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TIC Integration Parameters: LSCINT.P

Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.075	2.93 ng	357705	1,4-Dichlorobenzene-d4	7.934

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	78
2			Pentane, 3-methoxy-	102	C6H14O	036839-67-5	23
3			Silane, tetramethyl-	88	C4H12Si	000075-76-3	17
4			N-Ethylformamide	73	C3H7NO	000627-45-2	9
5			1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	9



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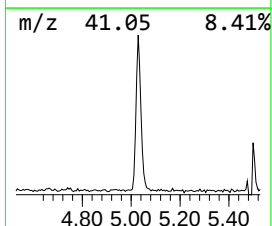
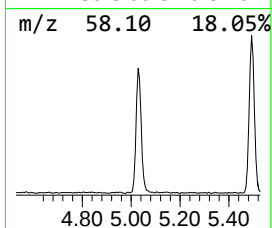
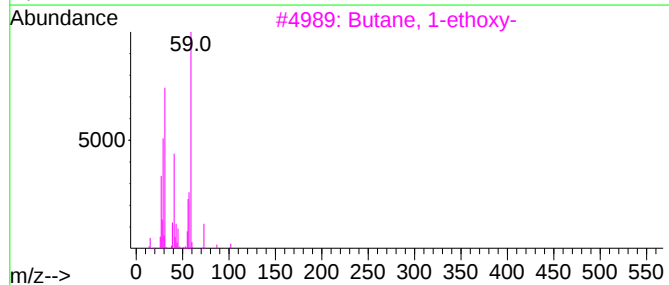
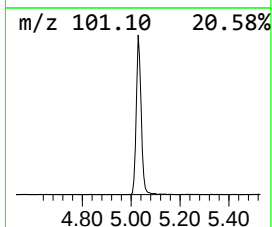
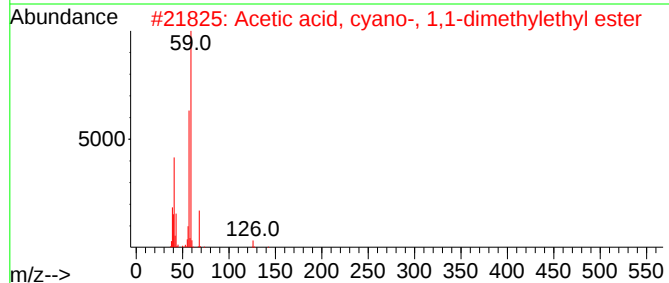
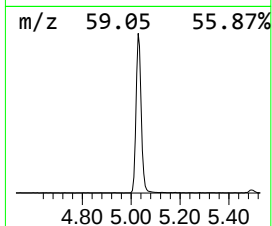
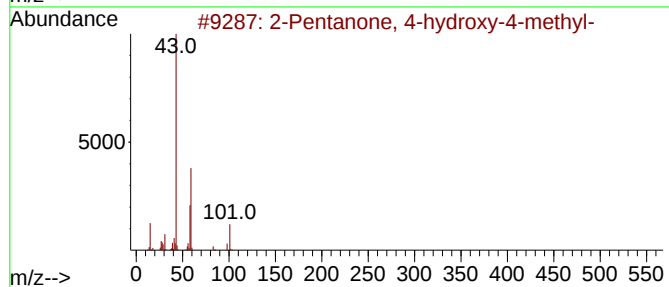
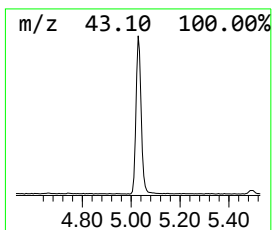
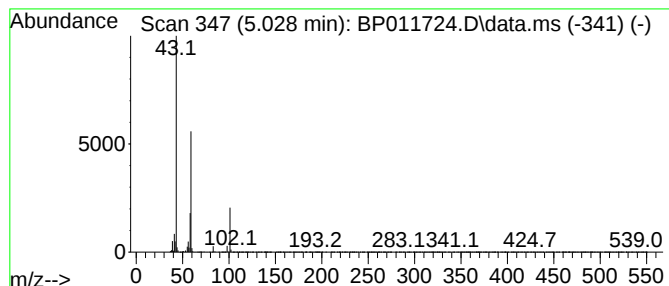
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Peak Number 2 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.028	6.77 ng	826631	1,4-Dichlorobenzene-d4	7.934

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50
2			Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	17
3			Butane, 1-ethoxy-	102	C6H14O	000628-81-9	9
4			Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9
5			Propane, 1-ethoxy-2-methyl-	102	C6H14O	000627-02-1	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Butane, 2-metho...	3.075	2.9	ng	357705	1	7.934	2442040	20.0
2-Pentanone, 4-...	5.028	6.8	ng	826631	1	7.934	2442040	20.0