

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042225\
 Data File : BP024374.D
 Acq On : 22 Apr 2025 11:54
 Operator : RC/JU
 Sample : PB167670BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 PB167670BL

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-BP041425.M
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BP024374.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	4.887	300	306	320	rBV	970075	1316190	8.52%	1.824%
2	5.346	377	384	404	rBV	5221645	7536279	48.78%	10.446%
3	5.940	479	485	494	rBV	193364	268230	1.74%	0.372%
4	6.911	640	650	681	rBV	4828662	7449446	48.21%	10.326%
5	7.716	781	787	797	rBV	953442	1510571	9.78%	2.094%
6	8.869	975	983	999	rBV	2924630	4839346	31.32%	6.708%
7	10.493	1250	1259	1271	rBV	1147300	2062392	13.35%	2.859%
8	12.957	1670	1678	1696	rBV	7610338	10964175	70.96%	15.198%
9	14.345	1907	1914	1925	rBV	1768094	2614196	16.92%	3.624%
10	15.863	2163	2172	2195	rBV	5768400	8699907	56.31%	12.059%
11	17.145	2383	2390	2403	rBV	1992224	2952880	19.11%	4.093%
12	19.869	2846	2853	2866	rBV	12676460	15450946	100.00%	21.417%
13	21.580	3138	3144	3157	rBV2	1985163	3226831	20.88%	4.473%
14	24.927	3703	3713	3727	rVB	1050543	3252575	21.05%	4.508%

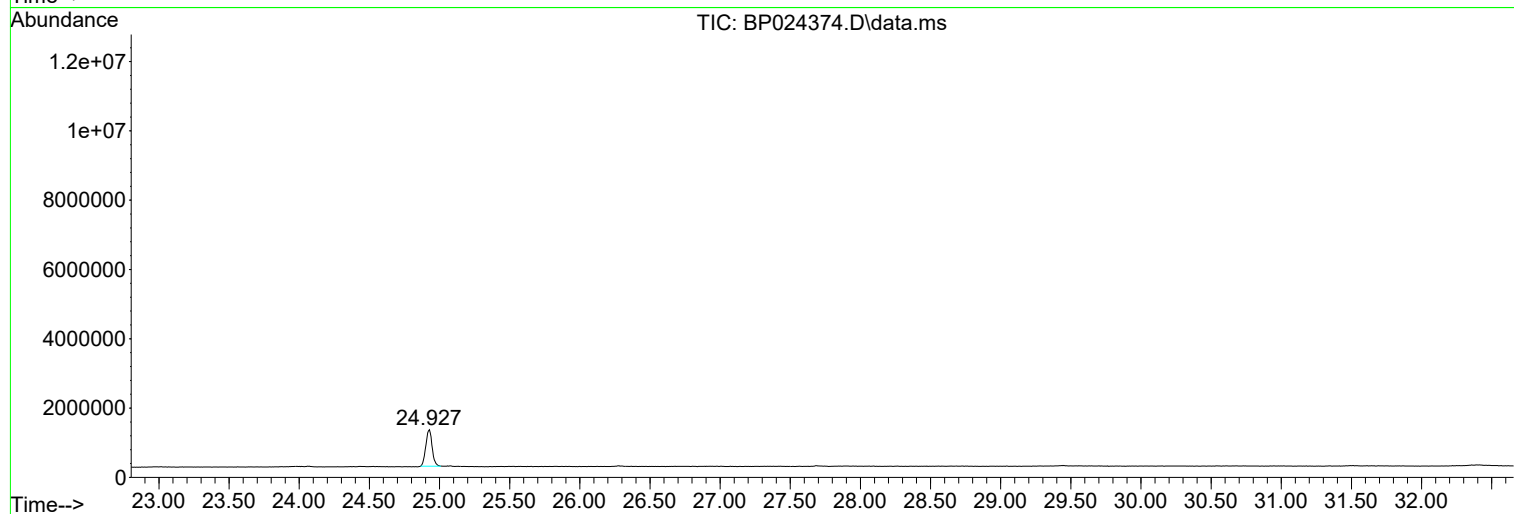
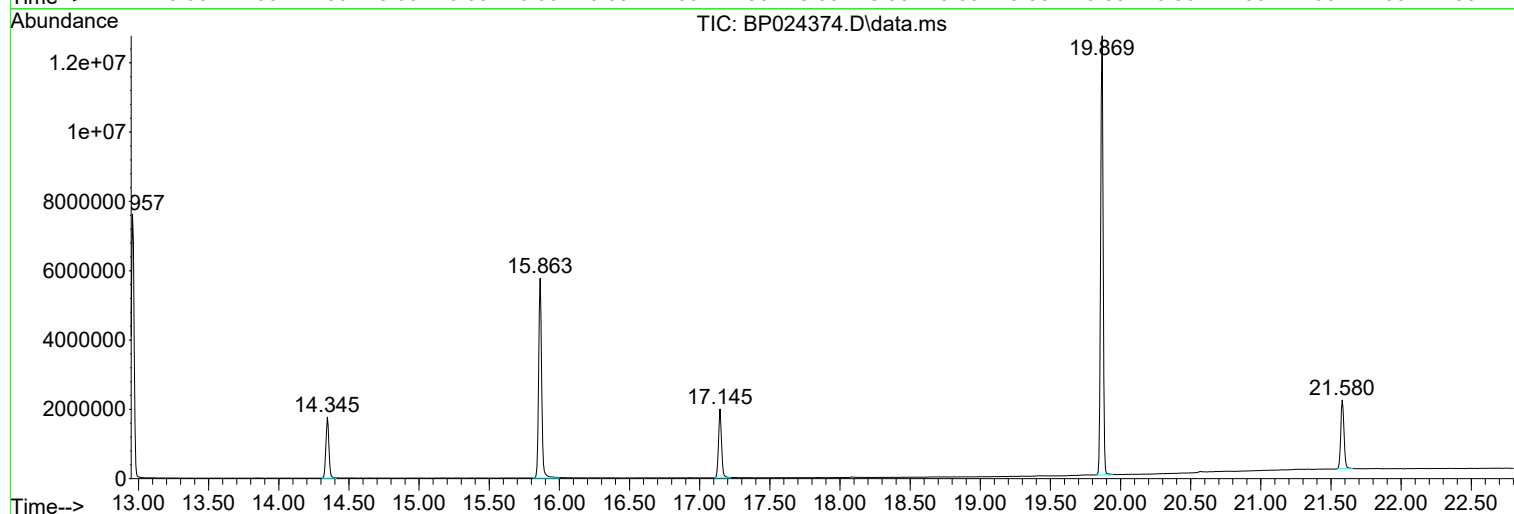
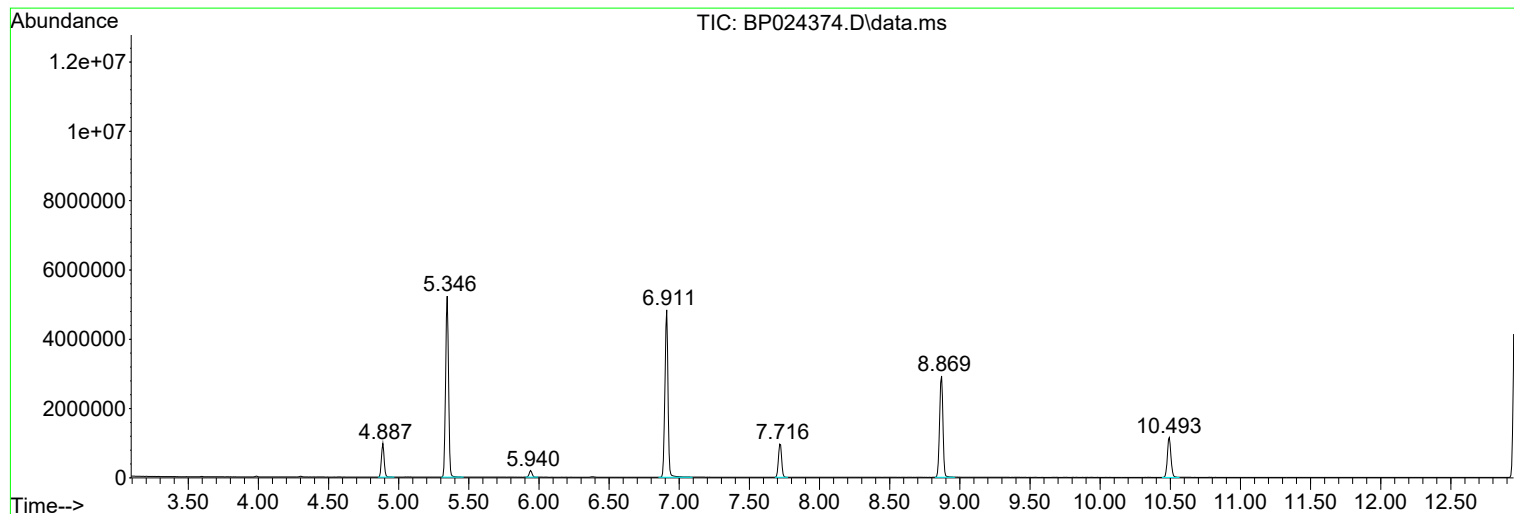
Sum of corrected areas: 72143964

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP041425.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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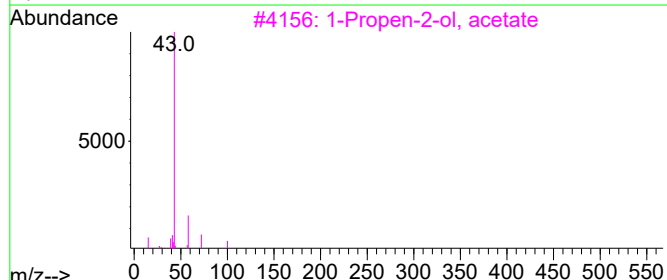
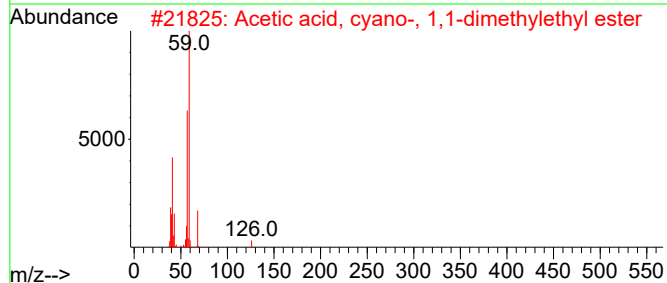
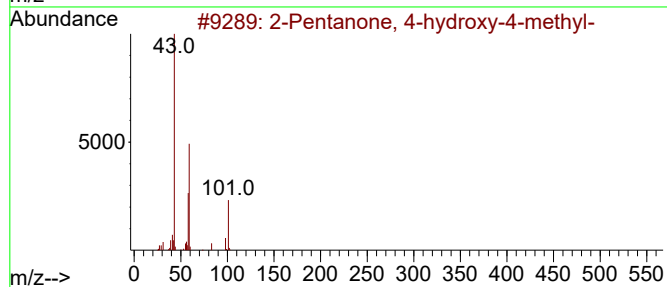
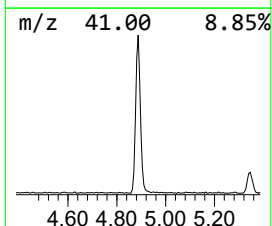
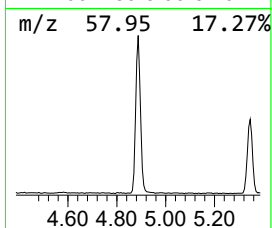
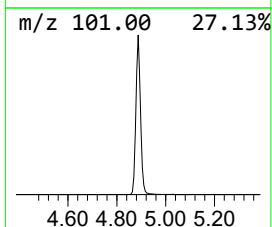
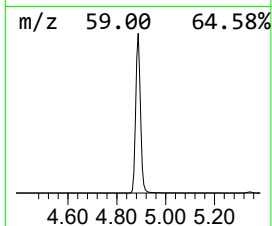
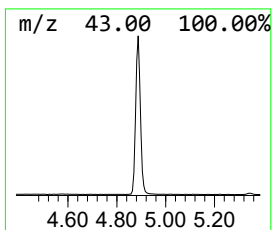
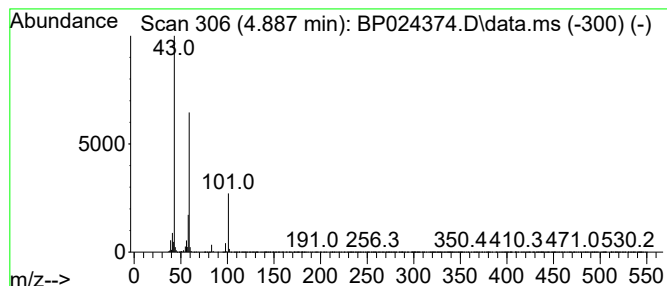
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TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD			R.T.
4.887	17.43 ng	1316190	1,4-Dichlorobenzene-d4			7.722
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-		116	C6H12O2	000123-42-2	72
2	Acetic acid, cyano-, 1,1-dimethy...		141	C7H11NO2	001116-98-9	17
3	1-Propen-2-ol, acetate		100	C5H8O2	000108-22-5	10
4	5-Hexen-2-one		98	C6H10O	000109-49-9	9
5	2,3-Butanedione, monooxime		101	C4H7NO2	000057-71-6	9



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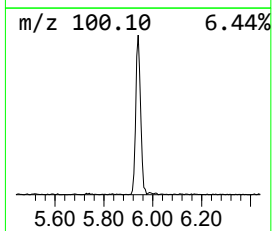
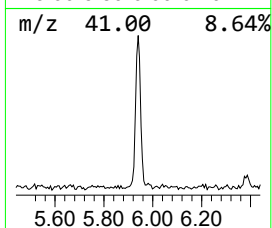
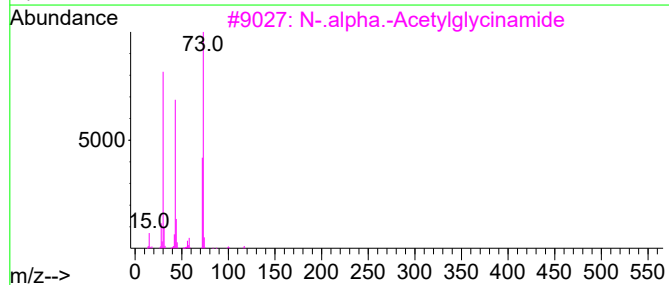
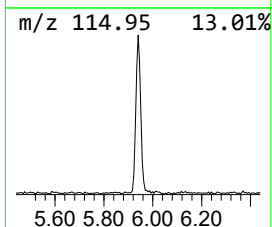
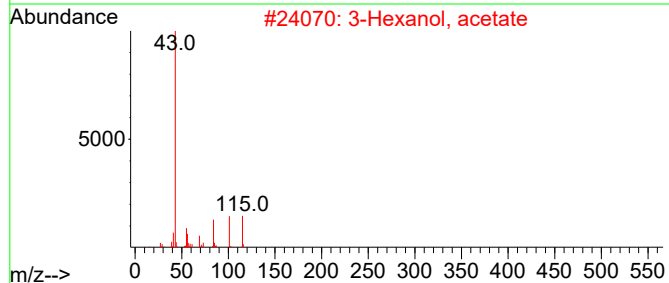
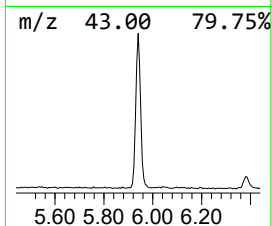
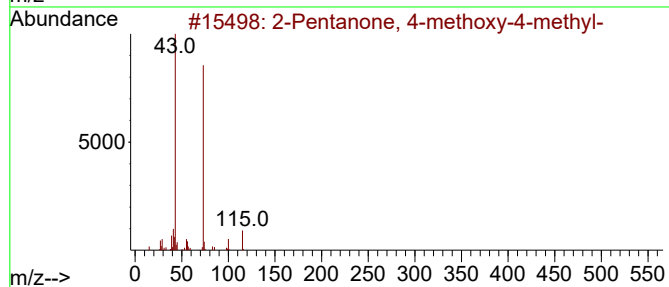
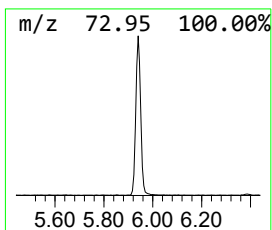
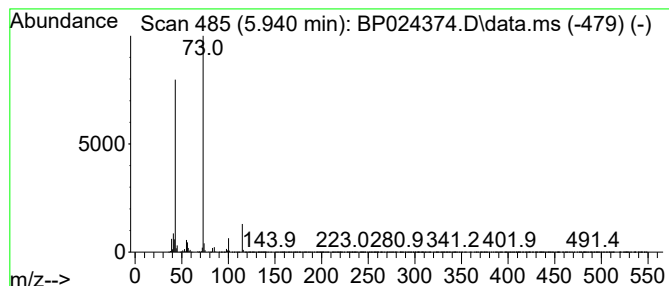
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Peak Number 2 2-Pentanone, 4-methoxy-4-me... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.940	3.55 ng	268230	1,4-Dichlorobenzene-d4	7.722

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-methoxy-4-methyl-	130	C7H14O2	000107-70-0	83
2			3-Hexanol, acetate	144	C8H16O2	040780-64-1	37
3			N-.alpha.-Acetylglycinamide	116	C4H8N2O2	002620-63-5	28
4			2-Pentanone, 3-methyl-	100	C6H12O	000565-61-7	22
5			2-Pentanone, 3-ethyl-	114	C7H14O	006137-03-7	10



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-...	4.887	17.4	ng	1316190	1	7.722	1510570	20.0
2-Pentanone, 4-...	5.940	3.5	ng	268230	1	7.722	1510570	20.0