

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041525\
 Data File : BP024287.D
 Acq On : 15 Apr 2025 09:56
 Operator : RC/JU
 Sample : PB167571BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 PB167571BL

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: BP024287.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	301	306	316	rBV	794486	1106835	8.32%	1.804%
2	5.346	378	384	406	rBV	4266766	6153200	46.25%	10.028%
3	5.946	481	486	496	rVB	138170	201534	1.51%	0.328%
4	6.911	642	650	670	rBV	3870684	6197806	46.59%	10.101%
5	7.728	782	789	796	rBV	874641	1302672	9.79%	2.123%
6	8.875	975	984	1005	rBV	2501986	3996124	30.04%	6.513%
7	10.505	1251	1261	1276	rBV	996554	1816312	13.65%	2.960%
8	12.969	1672	1680	1696	rBV	6272938	8941071	67.21%	14.572%
9	14.363	1910	1917	1930	rBV	1402443	2331807	17.53%	3.800%
10	15.881	2167	2175	2194	rBV	5333380	6814676	51.23%	11.106%
11	17.175	2388	2395	2409	rBV2	1753629	2652894	19.94%	4.324%
12	19.910	2854	2860	2870	rBV	10627867	13303149	100.00%	21.681%
13	21.616	3144	3150	3165	rBV	1687675	3151486	23.69%	5.136%
14	24.992	3714	3724	3741	rVB	1117720	3389497	25.48%	5.524%

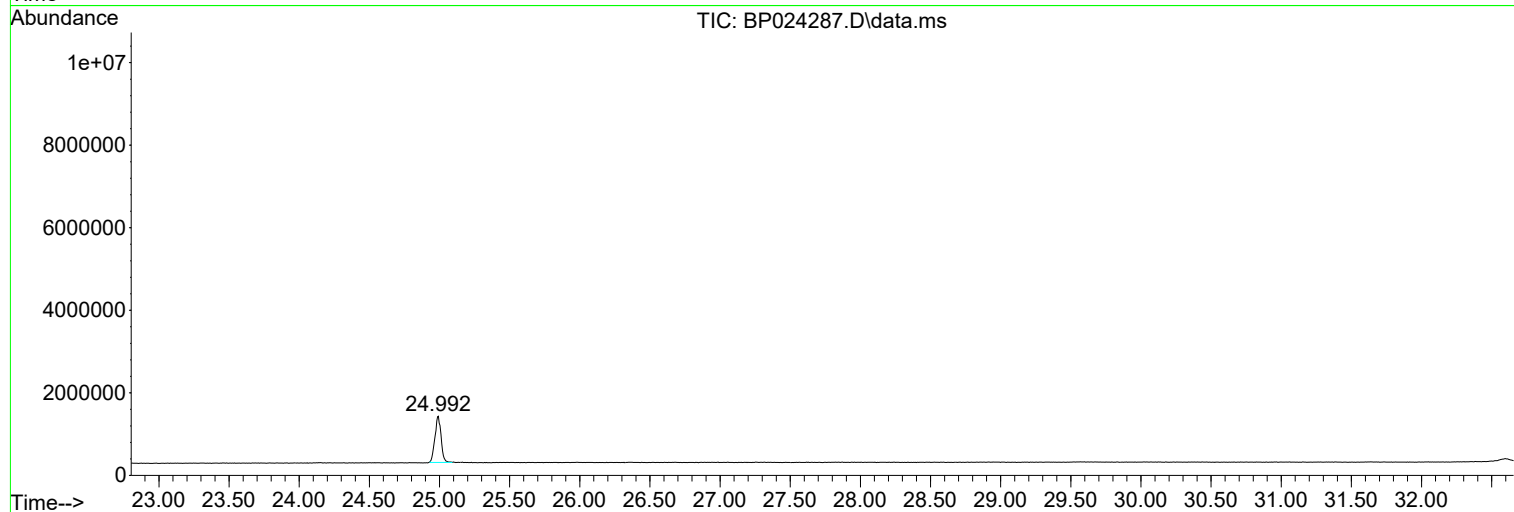
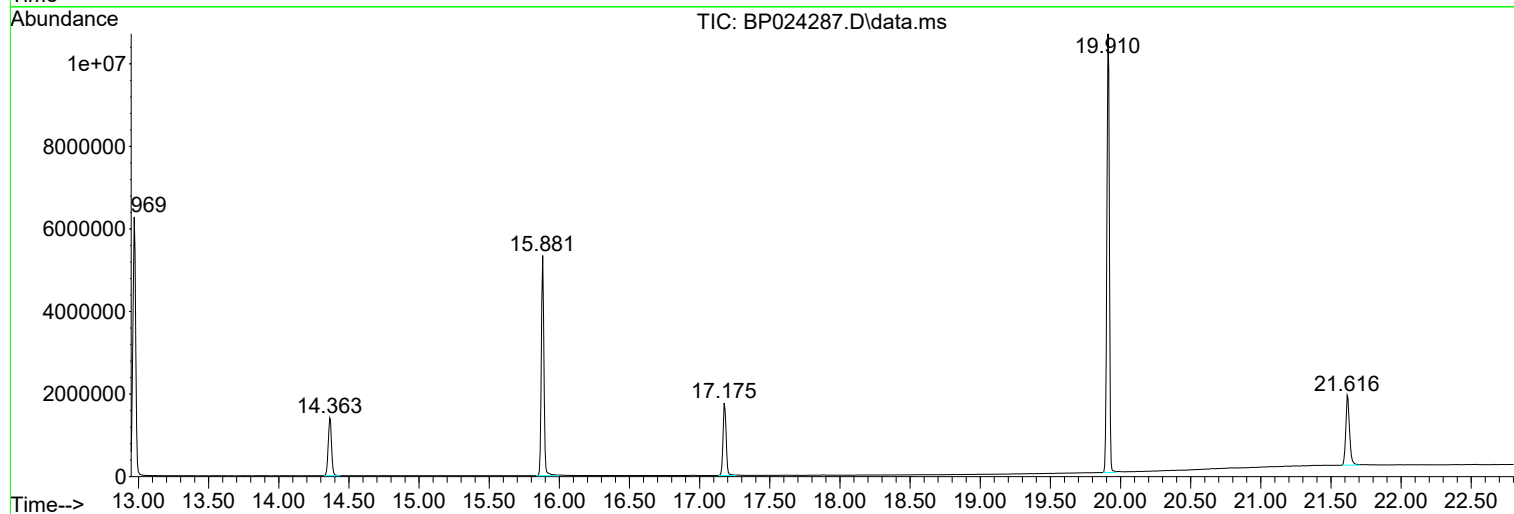
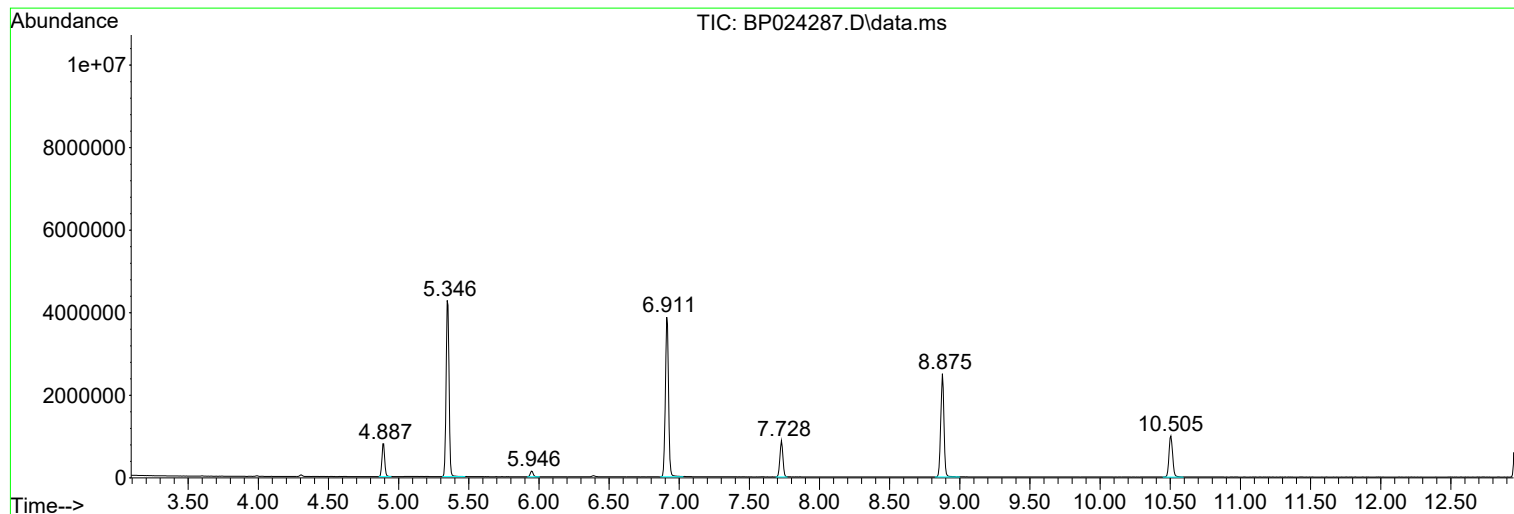
Sum of corrected areas: 61359063

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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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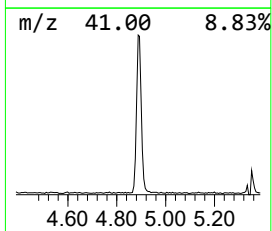
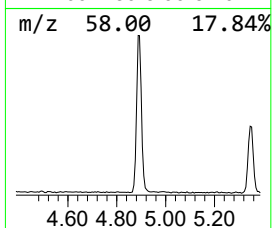
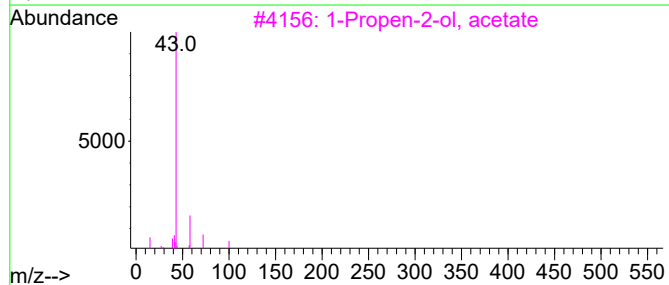
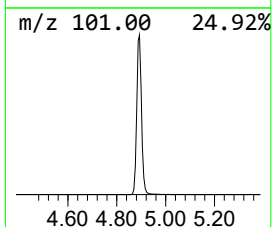
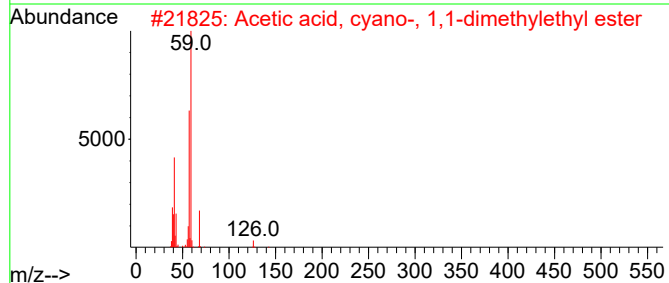
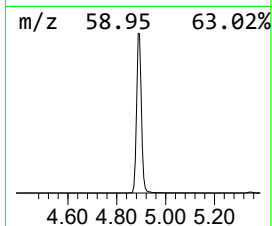
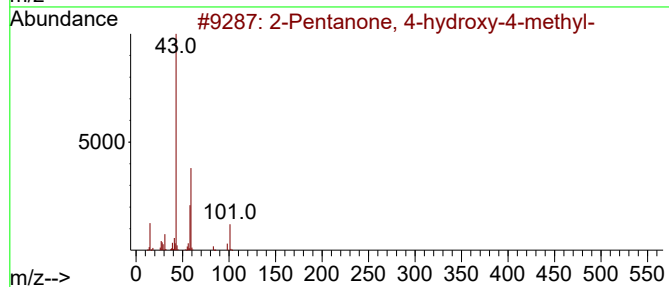
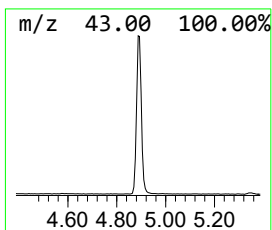
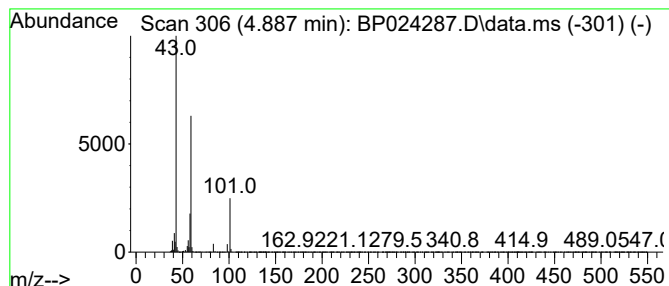
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.887	16.99 ng	1106840	1,4-Dichlorobenzene-d4	7.728

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	17		
3	1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	10		
4	Butane, 1-ethoxy-	102	C6H14O	000628-81-9	9		
5	(+)-4-Amino-4,5-dihydro-2(3H)-f...	101	C4H7NO2	016504-58-8	9		



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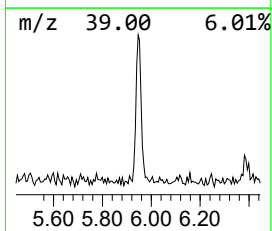
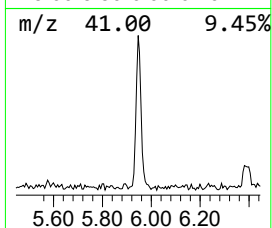
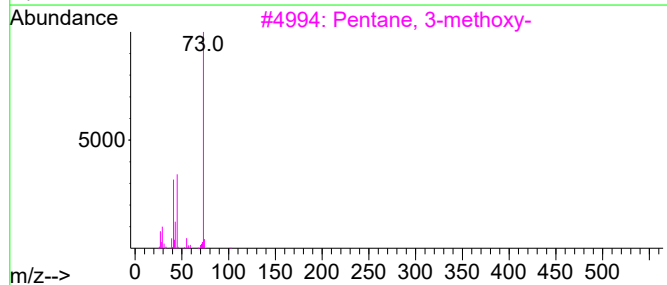
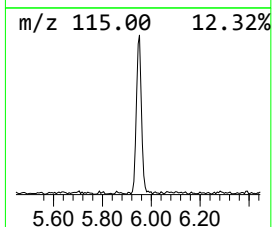
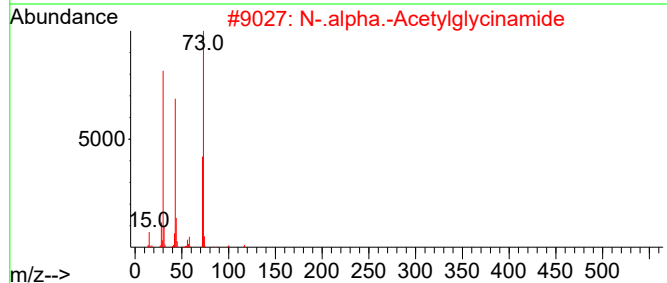
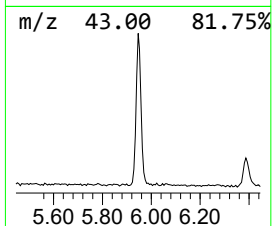
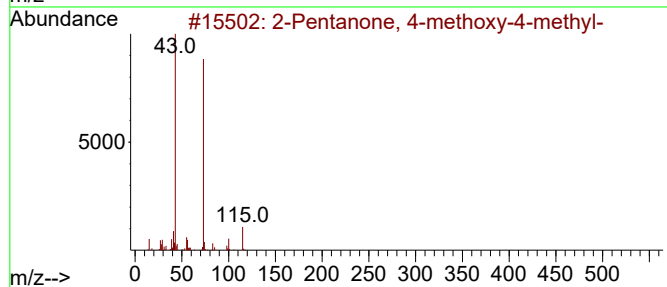
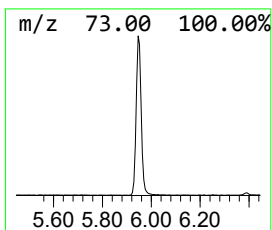
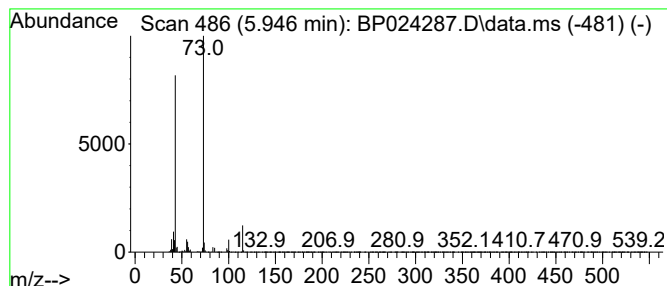
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Peak Number 2 2-Pentanone, 4-methoxy-4-me... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.946	3.09 ng	201534	1,4-Dichlorobenzene-d4	7.728

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		2-Pentanone, 4-methoxy-4-methyl-	130	C7H14O2	000107-70-0	83
2		N-.alpha.-Acetylglycinamide	116	C4H8N2O2	002620-63-5	33
3		Pentane, 3-methoxy-	102	C6H14O	036839-67-5	12
4		4-Penten-2-one, 4-methyl-	98	C6H10O	003744-02-3	10
5		1,3-Dioxolane, 2-(1-methylethyl)-	116	C6H12O2	000822-83-3	10



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TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-...	4.887	17.0	ng	1106840	1	7.728	1302670	20.0
2-Pentanone, 4-...	5.946	3.1	ng	201534	1	7.728	1302670	20.0