

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041625\
 Data File : BP024309.D
 Acq On : 16 Apr 2025 13:07
 Operator : RC/JU
 Sample : PB167564BL
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 PB167564BL

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: BP024309.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	4.893	301	307	322	rBV	1391546	1925943	13.10%	2.697%
2	5.346	378	384	399	rBV	5094960	7194349	48.95%	10.076%
3	5.945	480	486	498	rVB	270849	382631	2.60%	0.536%
4	6.910	641	650	673	rBV	4541677	7245027	49.30%	10.147%
5	7.728	781	789	801	rBV	1010101	1530529	10.41%	2.144%
6	8.875	976	984	1001	rBV	2840104	4693233	31.93%	6.573%
7	10.504	1251	1261	1280	rBV	1250394	2128003	14.48%	2.980%
8	12.969	1672	1680	1698	rBV	6879725	10929755	74.37%	15.308%
9	14.363	1909	1917	1928	rBV	1797662	2750664	18.72%	3.853%
10	15.874	2167	2174	2206	rBV	5647508	8225069	55.97%	11.520%
11	17.174	2388	2395	2410	rBV	2254278	3043149	20.71%	4.262%
12	19.898	2853	2858	2874	rBV	11388528	14696517	100.00%	20.584%
13	21.627	3145	3152	3164	rBV2	1950638	3229365	21.97%	4.523%
14	24.992	3716	3724	3739	rVB	1331160	3424907	23.30%	4.797%

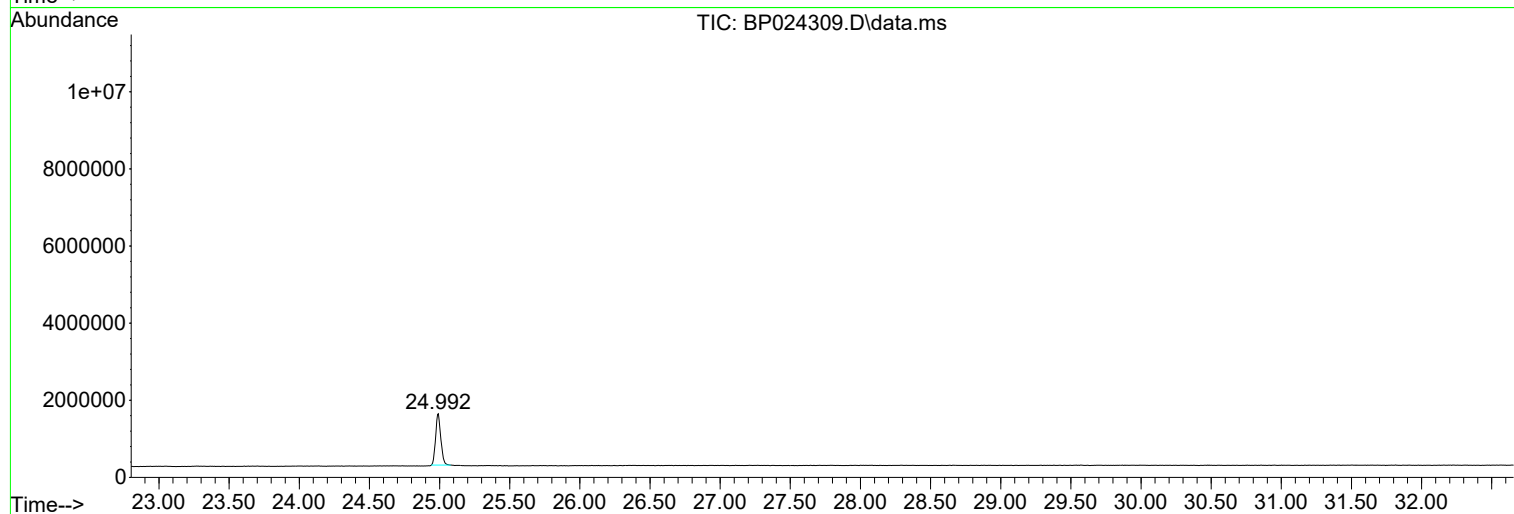
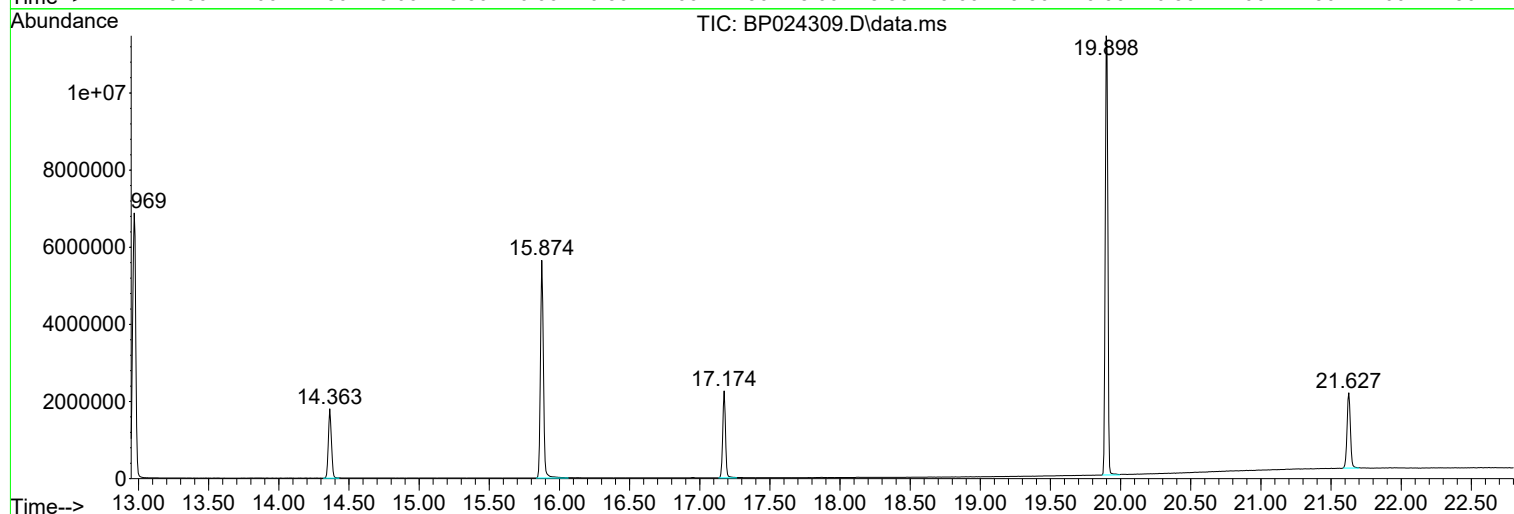
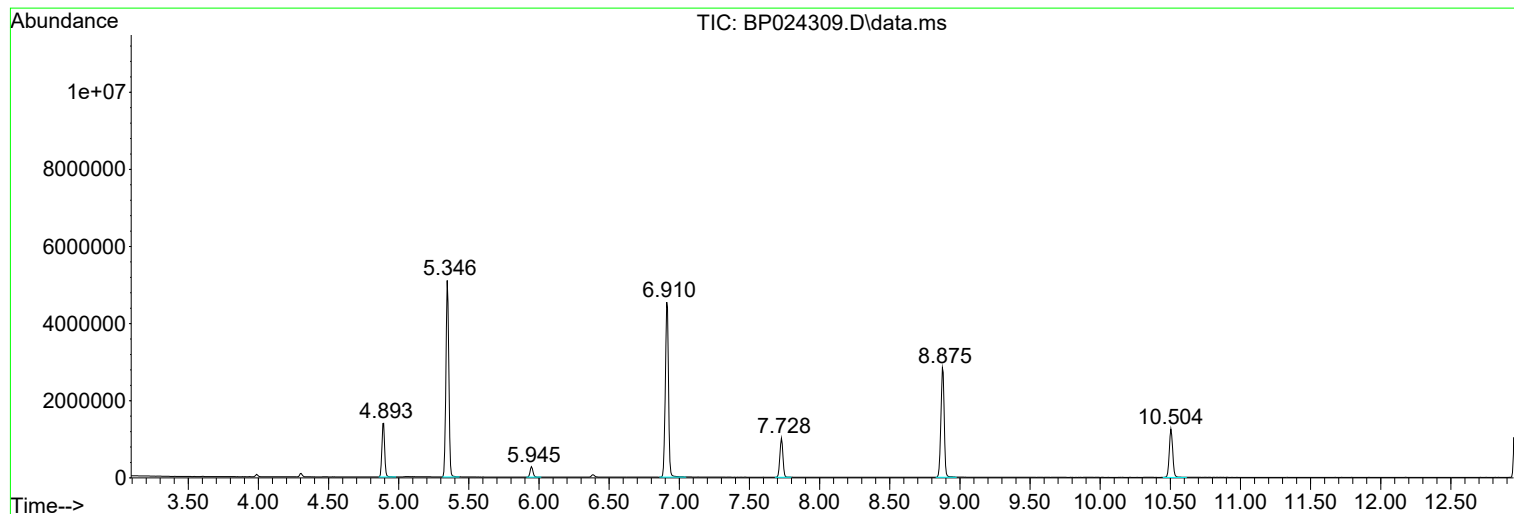
Sum of corrected areas: 71399141

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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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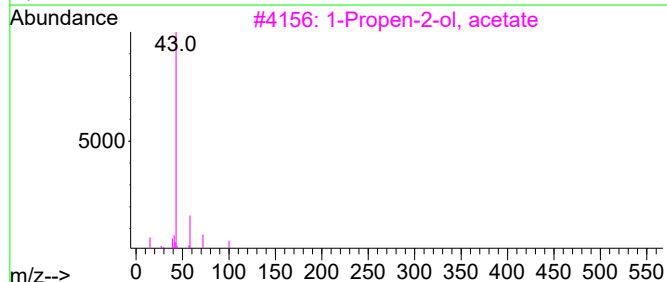
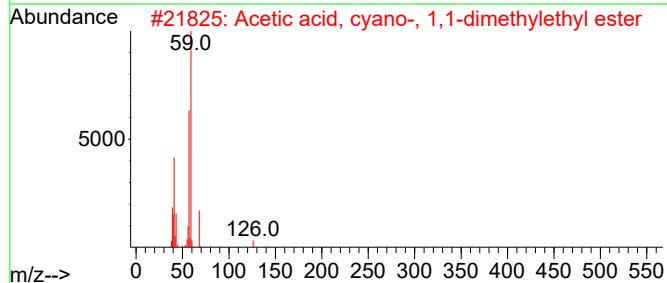
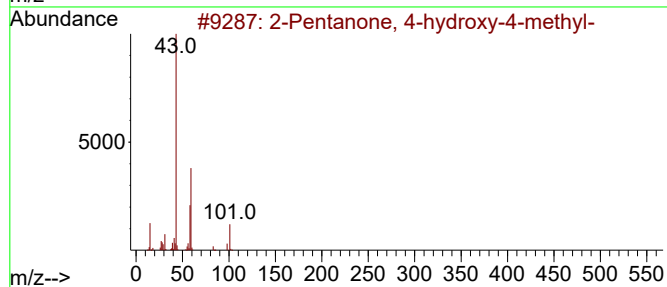
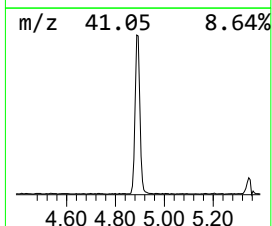
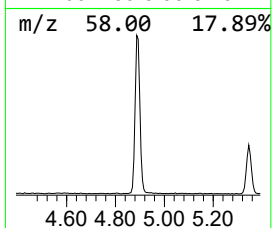
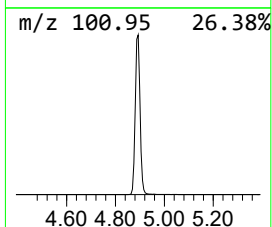
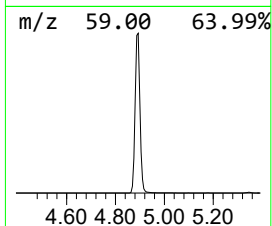
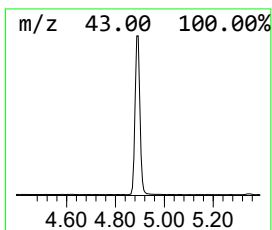
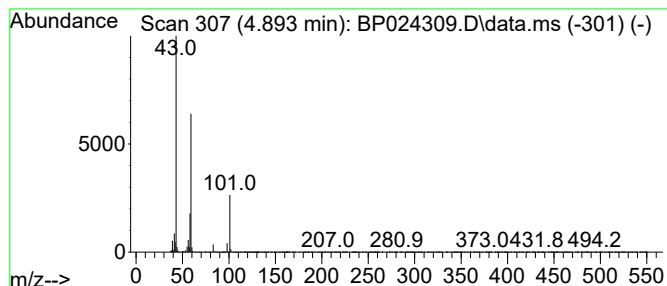
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TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.	
4.893	25.17 ng	1925940	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	5-Hexen-2-one	98	C6H10O	000109-49-9	9
5	Guanidine	59	CH5N3	000113-00-8	9



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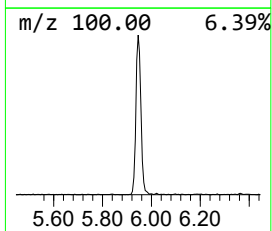
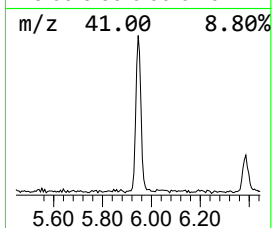
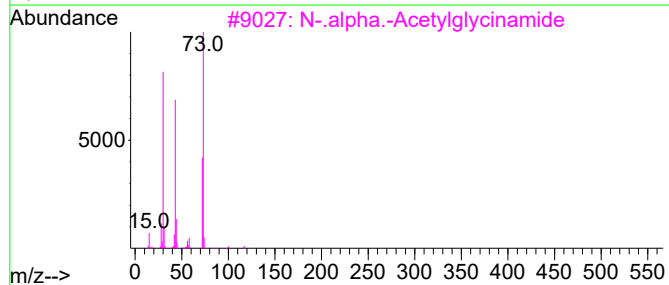
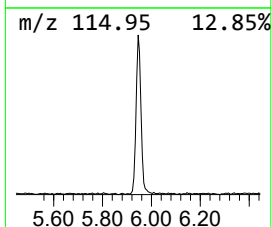
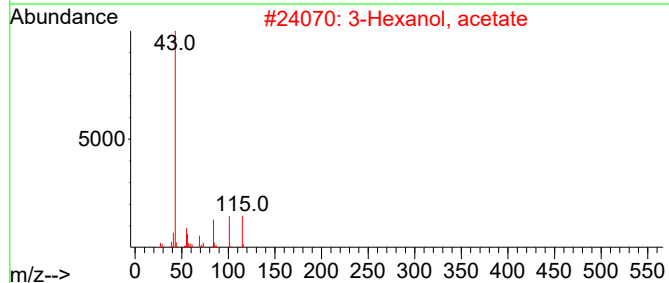
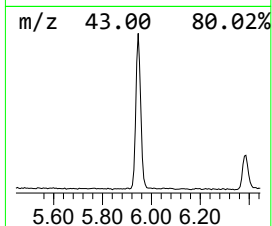
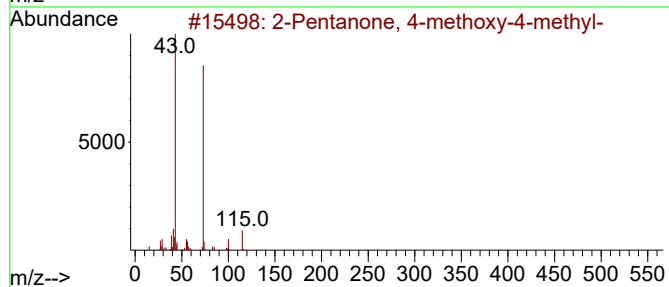
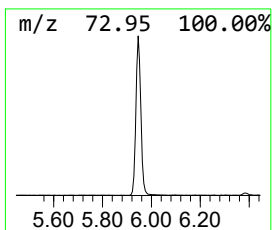
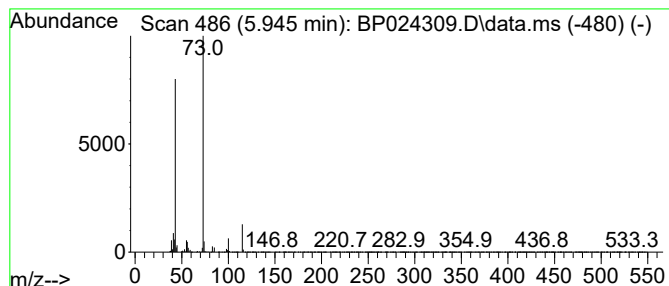
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Peak Number 2 2-Pentanone, 4-methoxy-4-me... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.945	5.00 ng	382631	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			3-Hexanol, acetate	144	C8H16O2	040780-64-1	37
3			N-.alpha.-Acetylglycinamide	116	C4H8N2O2	002620-63-5	28
4			Butane, 2-methoxy-2,3,3-trimethyl-	130	C8H18O	027705-21-1	23
5			Pentane, 3-methoxy-	102	C6H14O	036839-67-5	12



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-...	4.893	25.2	ng	1925940	1	7.728	1530530	20.0
2-Pentanone, 4-...	5.945	5.0	ng	382631	1	7.728	1530530	20.0