

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042425\
 Data File : BP024408.D
 Acq On : 24 Apr 2025 12:31
 Operator : RC/JU
 Sample : PB167711BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 PB167711BL

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: BP024408.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	320	rBV	1070085	1484539	10.52%	2.066%
2	5.346	377	384	398	rBV	5395189	7586279	53.75%	10.559%
3	5.946	480	486	496	rVB	223541	323061	2.29%	0.450%
4	6.910	641	650	671	rBV	4758018	7591408	53.78%	10.566%
5	7.722	780	788	799	rBV	1021605	1650819	11.70%	2.298%
6	8.869	975	983	1003	rBV	3066322	5043883	35.73%	7.020%
7	10.487	1249	1258	1275	rBV	1287312	2228990	15.79%	3.102%
8	12.957	1671	1678	1700	rBV	7922016	11423607	80.93%	15.899%
9	14.339	1907	1913	1931	rBV	1792174	2801226	19.85%	3.899%
10	15.857	2164	2171	2190	rBV	6027689	8690270	61.57%	12.095%
11	17.133	2381	2388	2404	rBV2	1862875	2988563	21.17%	4.159%
12	19.857	2845	2851	2861	rBV	11496888	14114939	100.00%	19.645%
13	21.574	3137	3143	3160	rBV	1759304	2961150	20.98%	4.121%
14	24.915	3703	3711	3727	rVB	1018980	2960577	20.97%	4.121%

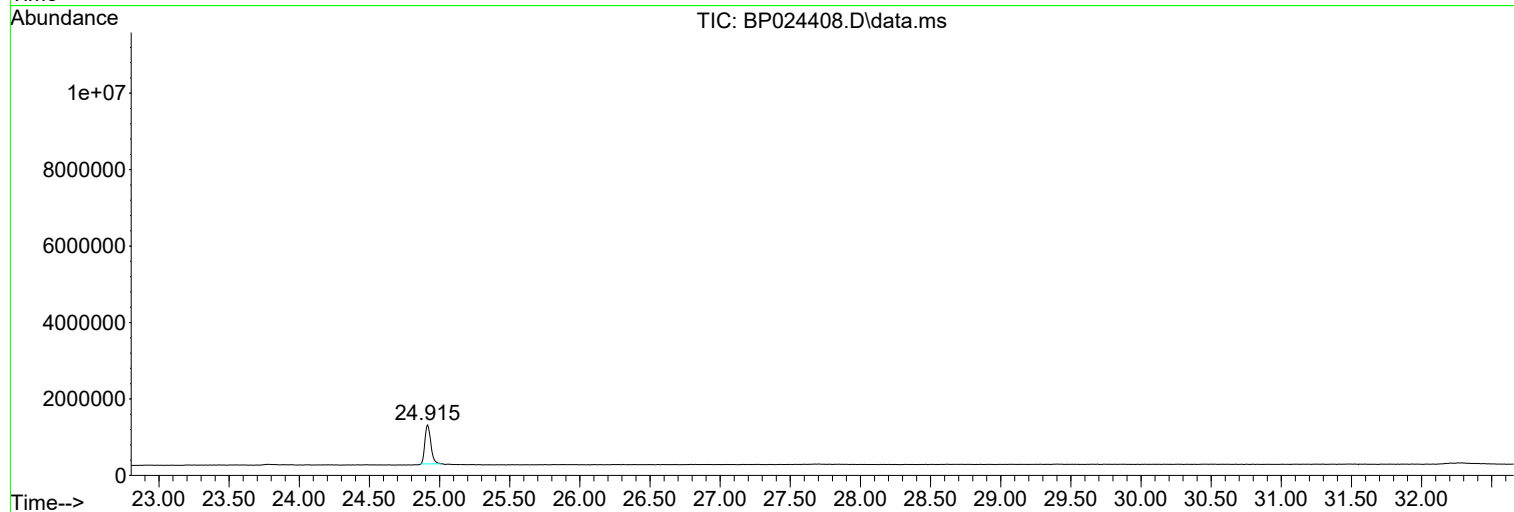
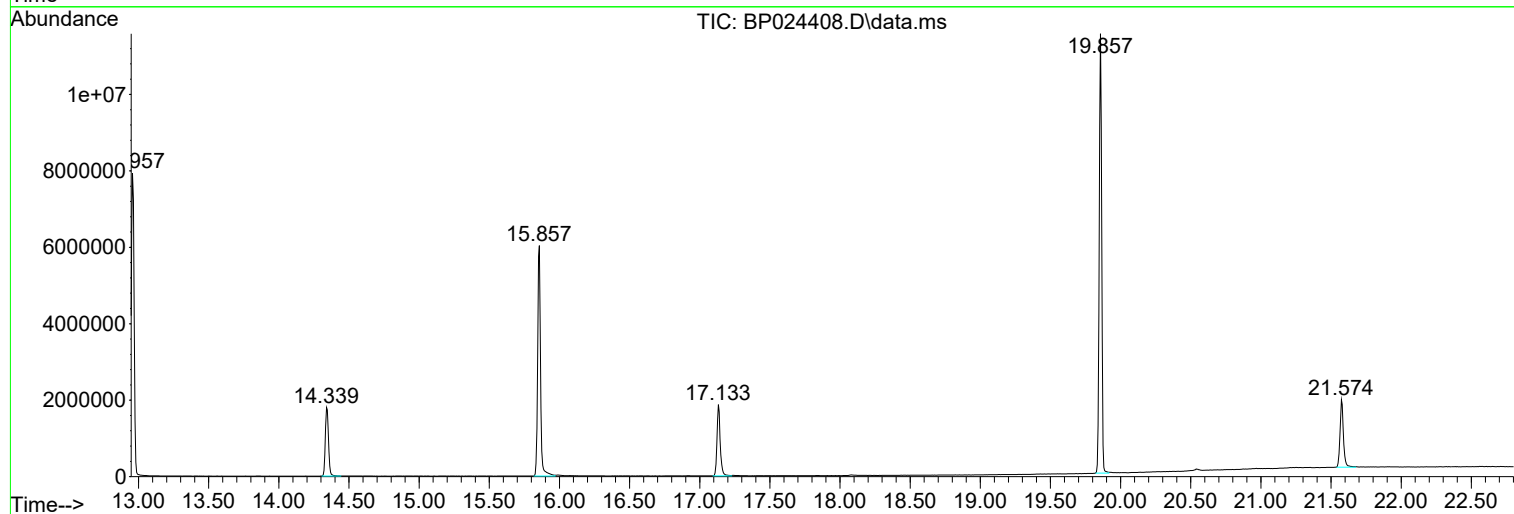
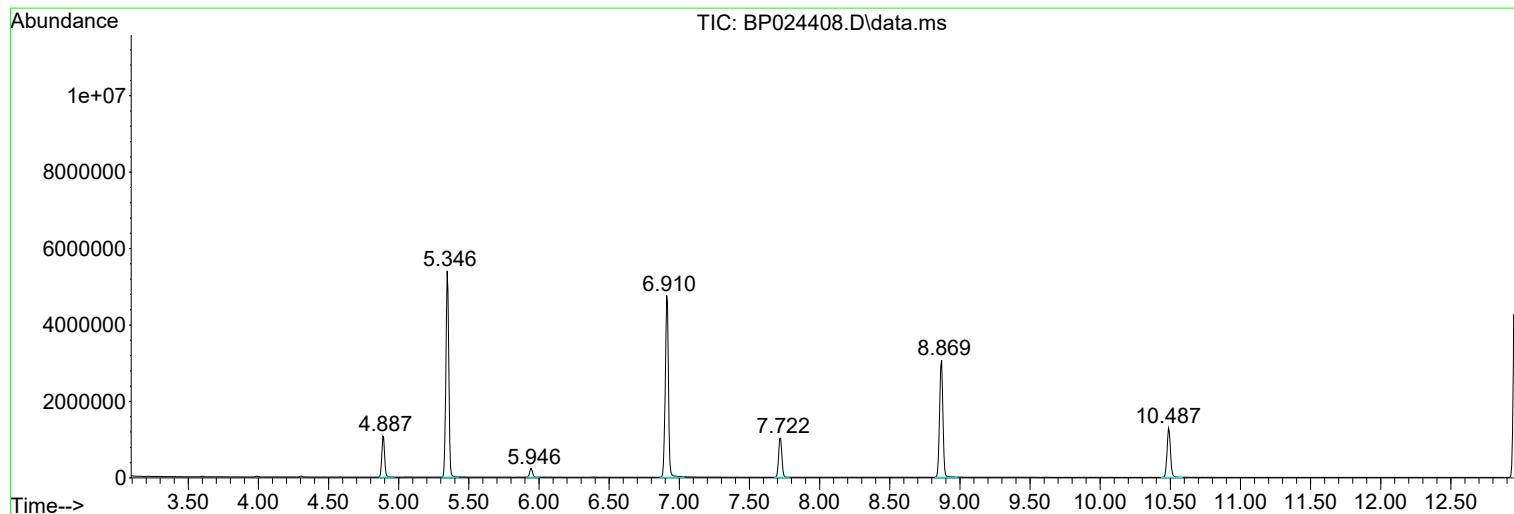
Sum of corrected areas: 71849311

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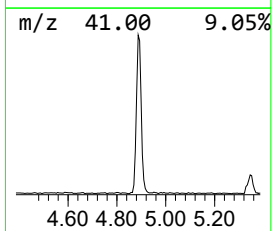
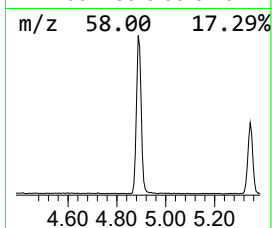
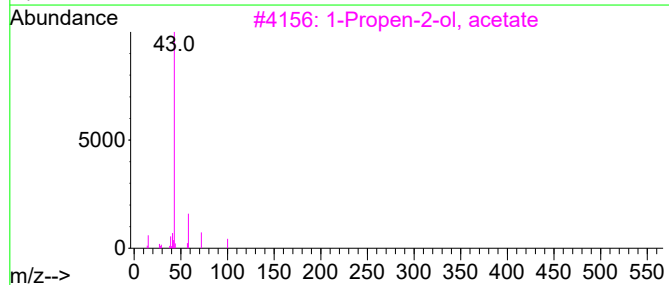
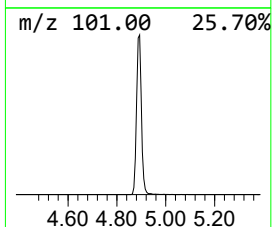
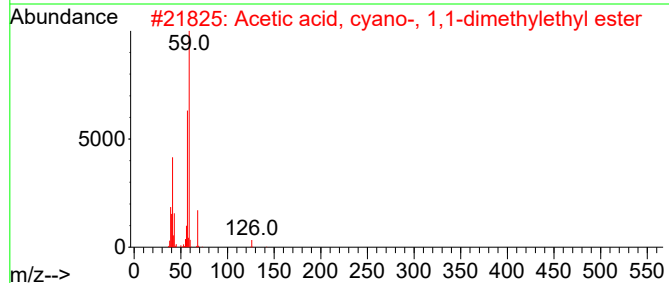
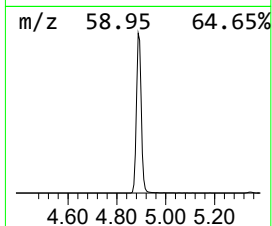
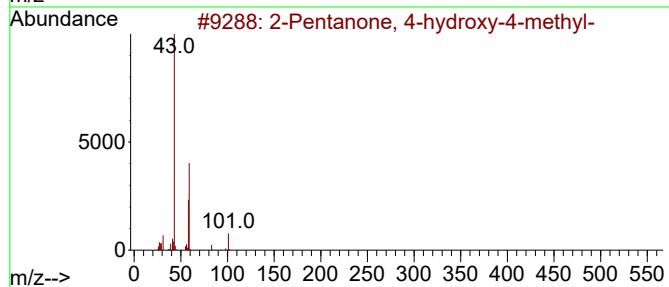
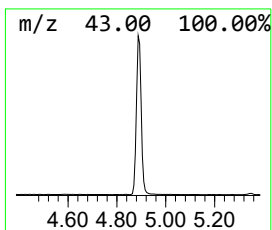
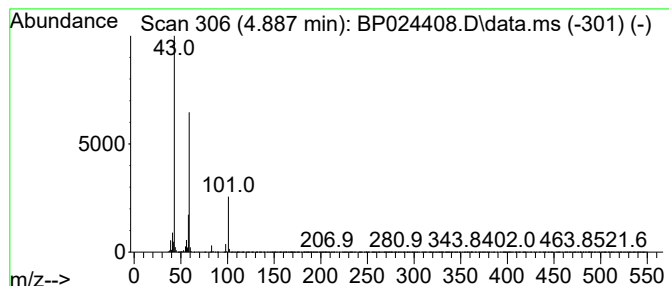
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.887	17.99 ng	1484540	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	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	Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9		
5	Butane, 1-ethoxy-	102	C6H14O	000628-81-9	9		



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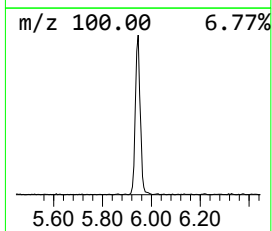
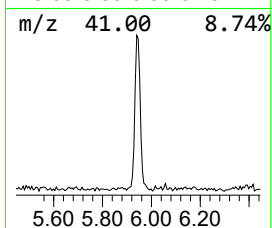
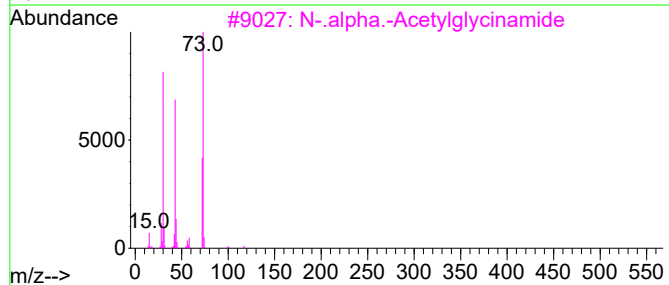
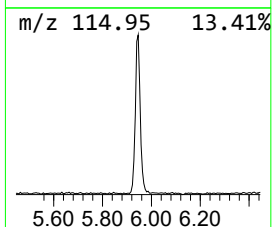
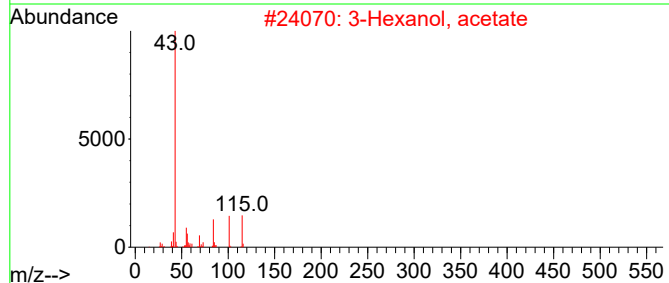
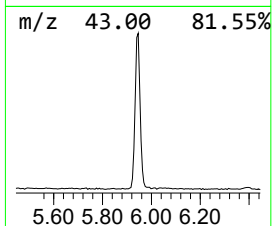
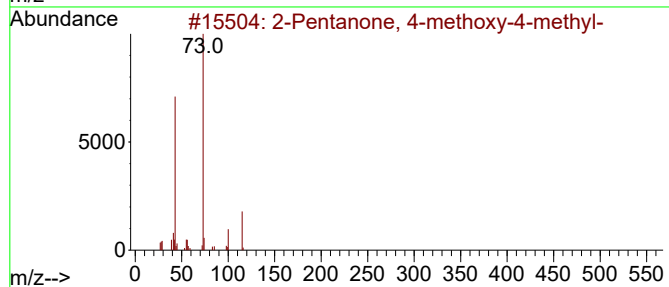
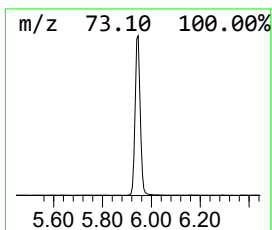
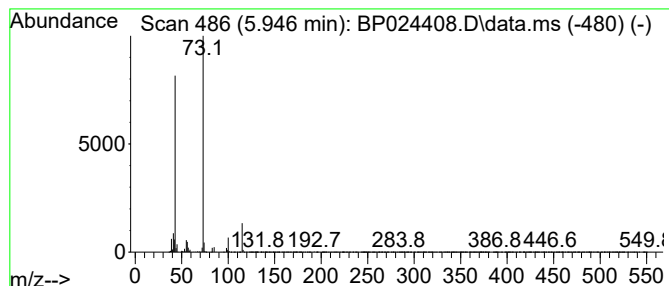
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Peak Number 2 2-Pentanone, 4-methoxy-4-me... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.946	3.91 ng	323061	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	32
3			N-.alpha.-Acetylglycinamide	116	C4H8N2O2	002620-63-5	28
4			2-Pentanone, 3-methyl-	100	C6H12O	000565-61-7	12
5			1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	12



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-...	4.887	18.0	ng	1484540	1	7.722	1650820	20.0
2-Pentanone, 4-...	5.946	3.9	ng	323061	1	7.722	1650820	20.0