

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF062525\
Data File : BF142838.D
Acq On : 25 Jun 2025 15:26
Operator : RC/JU
Sample : PB168601BL
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
PB168601BL

Integration Parameters: rteint.p

Integrator: RTE

Smoothing : ON

Sampling : 1

Start Thrs: 0.2

Stop Thrs : 0

Filtering: 5

Min Area: 3 % of largest Peak

Max Peaks: 100

Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >

Peak separation: 5

Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF061925.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BF142838.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.110	490	496	512	rBV	134562	218768	8.05%	1.393%
2	5.504	557	563	568	rBV	1567687	2094706	77.08%	13.340%
3	6.504	727	733	739	rBV	1415529	2031499	74.75%	12.937%
4	6.875	791	796	801	rBV	376505	456236	16.79%	2.905%
5	7.440	885	892	897	rBV	988208	1354913	49.86%	8.629%
6	8.157	1009	1014	1023	rBV	465796	594340	21.87%	3.785%
7	9.234	1191	1197	1203	rBV	1978735	2717613	100.00%	17.307%
8	9.916	1308	1313	1318	rBV	547186	673895	24.80%	4.292%
9	10.710	1442	1448	1462	rBV	1221589	1584076	58.29%	10.088%
10	11.404	1561	1566	1572	rBV	532159	669733	24.64%	4.265%
11	12.998	1831	1837	1842	rBV	1759942	2395333	88.14%	15.254%
12	14.051	2011	2016	2024	rBV	353759	447792	16.48%	2.852%
13	15.539	2263	2269	2279	rBV	294234	463780	17.07%	2.954%

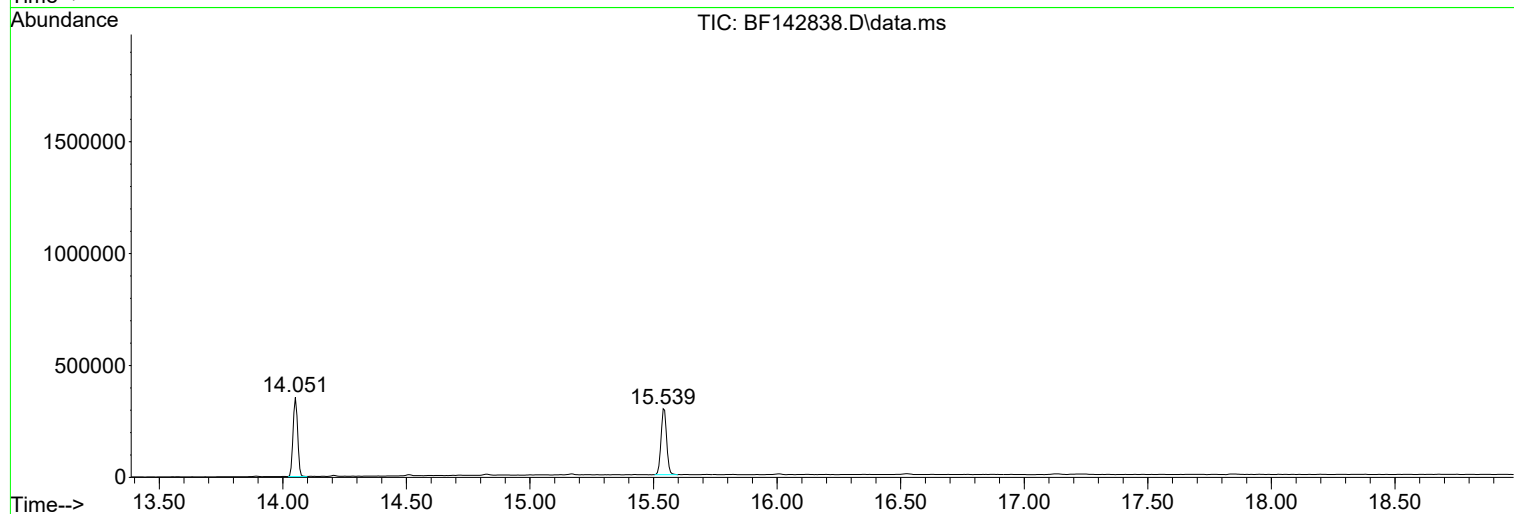
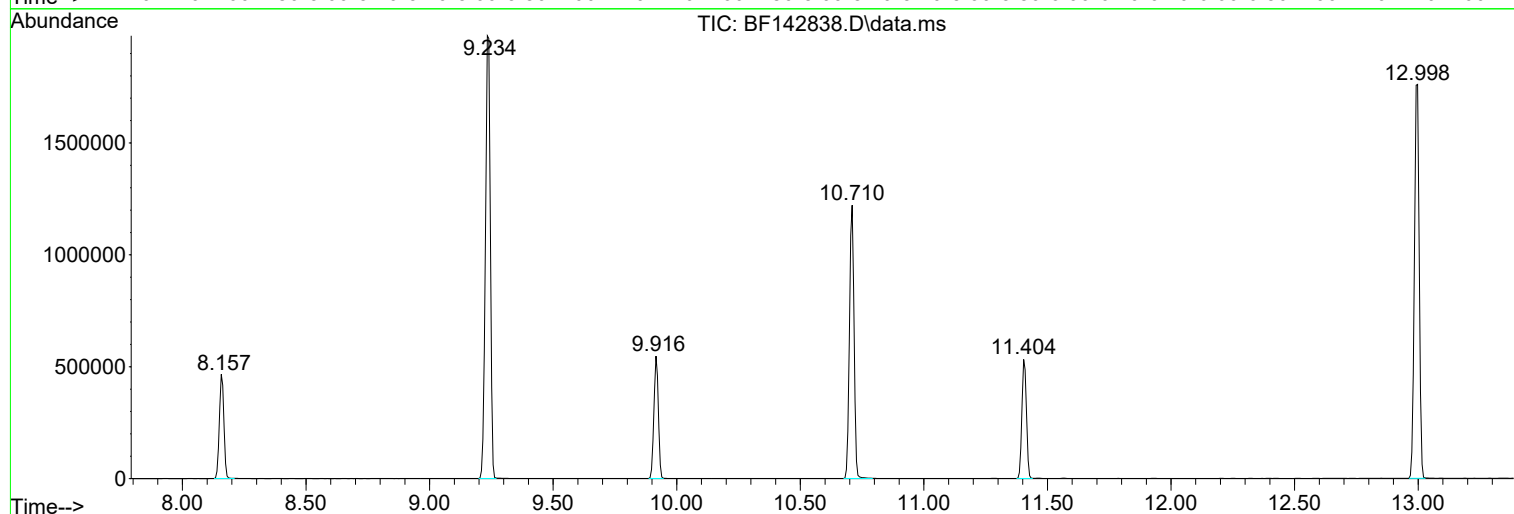
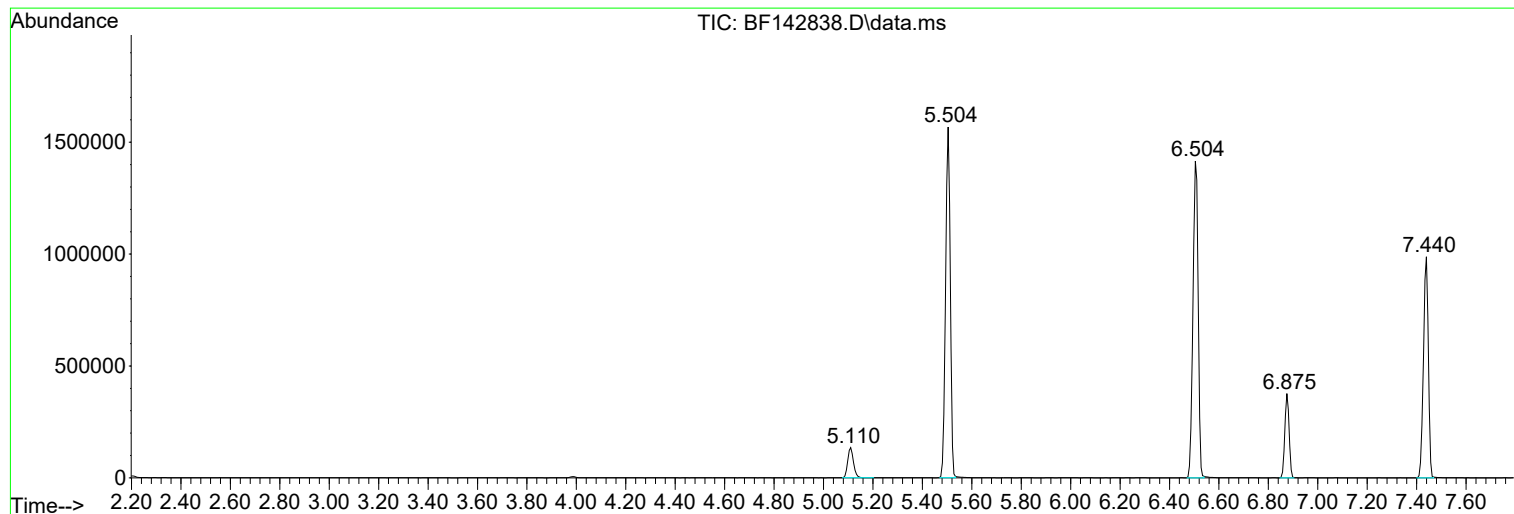
Sum of corrected areas: 15702684

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF061925.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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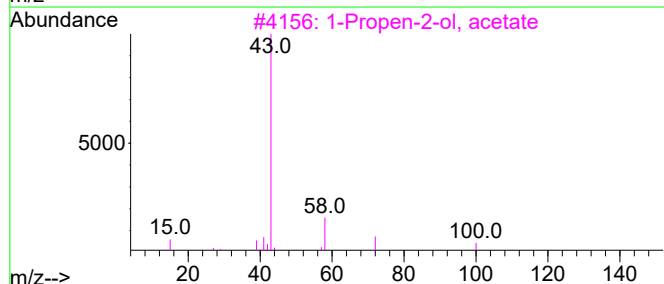
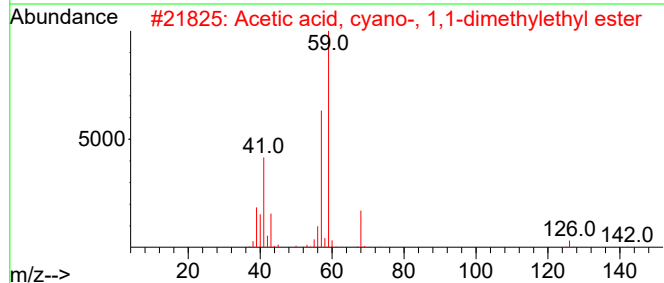
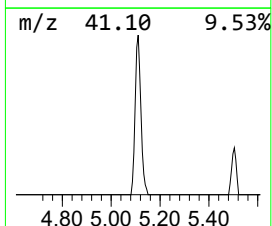
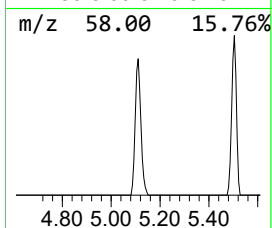
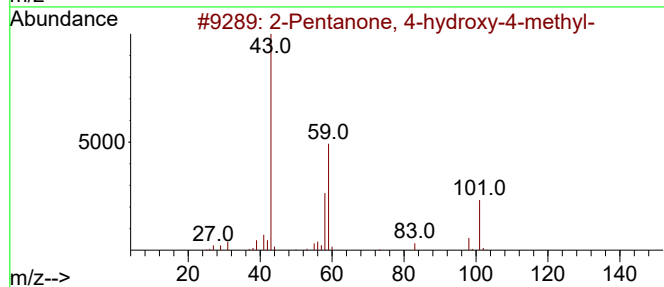
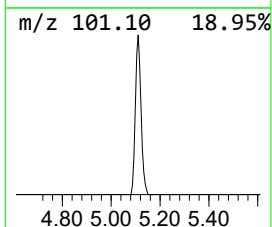
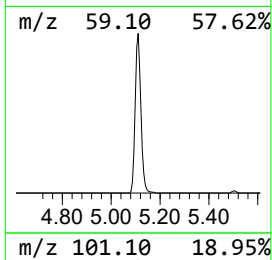
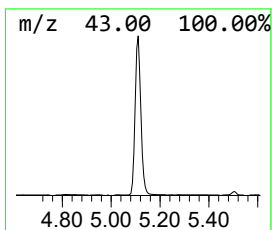
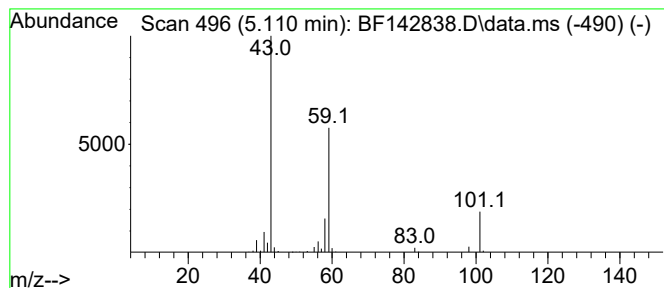
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.110	9.59 ng	218768	1,4-Dichlorobenzene-d4	6.875

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	32
3			1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	12
4			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	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
2-Pentanone, 4-...	5.110	9.6	ng	218768	1	6.875	456236	20.0