

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF060525\
Data File : BF142626.D
Acq On : 05 Jun 2025 11:04
Operator : RC/JU
Sample : PB168285BL
Misc :
ALS Vial : 3 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
PB168285BL

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_F\Methods\8270-BF052025.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BF142626.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.128	492	499	509	rBV	209595	338476	8.53%	1.465%
2	5.522	559	566	571	rBV	2271570	3086845	77.83%	13.358%
3	6.522	729	736	741	rBV	2128452	3043429	76.74%	13.170%
4	6.892	794	799	805	rBV	541774	681009	17.17%	2.947%
5	7.457	888	895	900	rBV	1480052	2025064	51.06%	8.764%
6	8.181	1012	1018	1033	rBV	681820	907366	22.88%	3.927%
7	9.257	1194	1201	1206	rBV	2963129	3965900	100.00%	17.163%
8	9.933	1311	1316	1327	rBV	817428	1042216	26.28%	4.510%
9	10.727	1445	1451	1471	rBV	1837266	2414649	60.89%	10.449%
10	11.421	1564	1569	1577	rBV	800851	1030686	25.99%	4.460%
11	13.010	1833	1839	1844	rBV	2549333	3343061	84.30%	14.467%
12	14.068	2011	2019	2033	rBV	434578	611688	15.42%	2.647%
13	15.562	2267	2273	2288	rVB	389081	617537	15.57%	2.672%

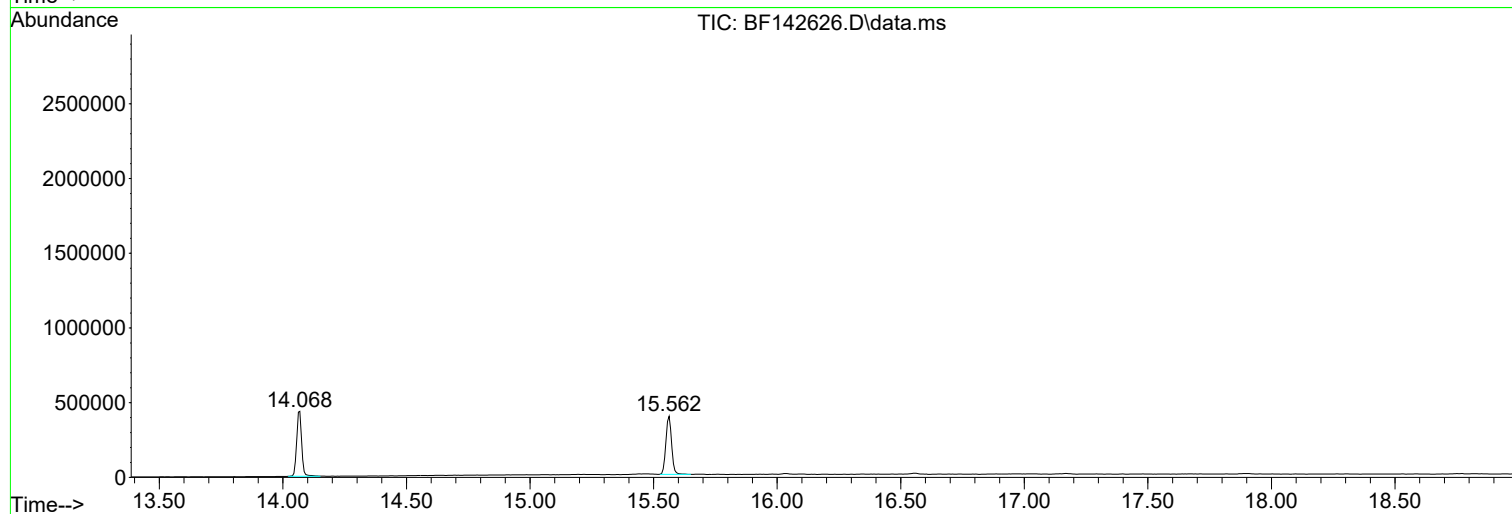
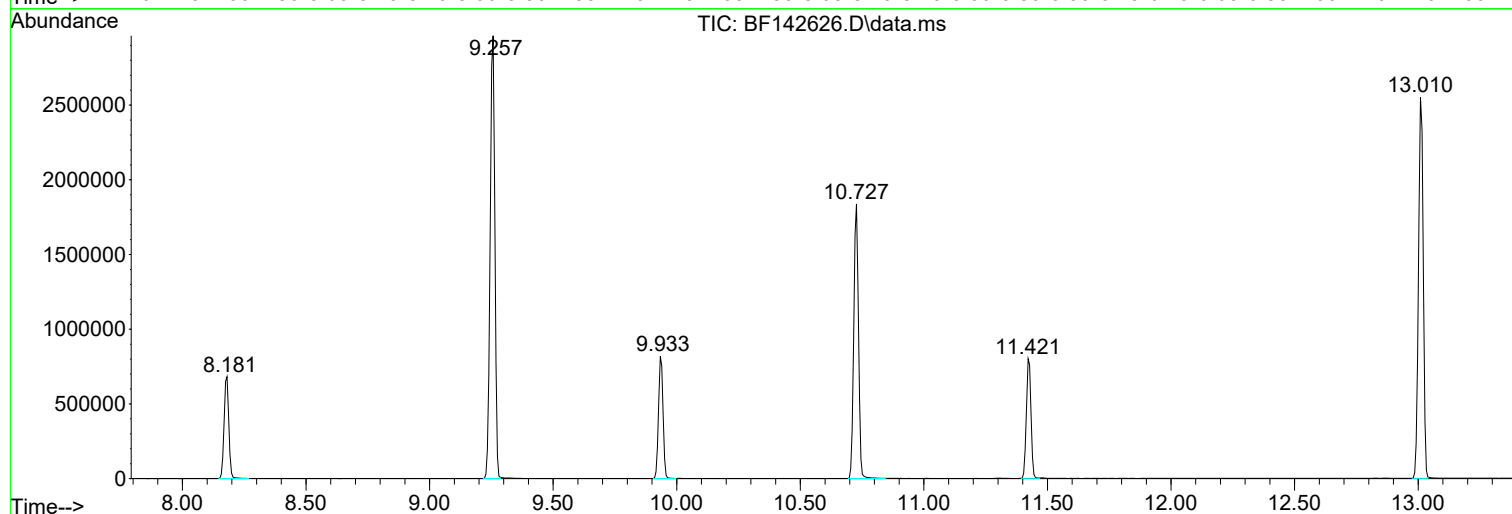
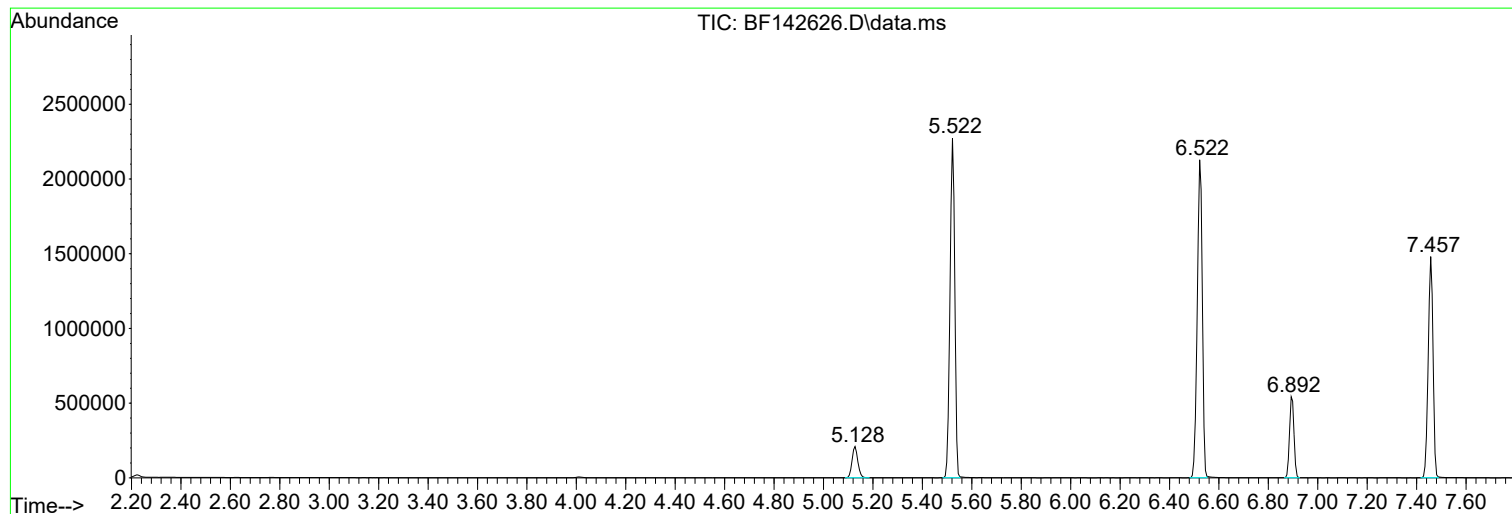
Sum of corrected areas: 23107926

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF052025.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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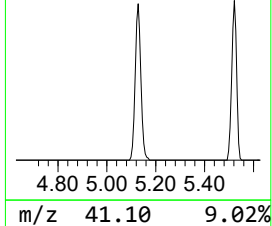
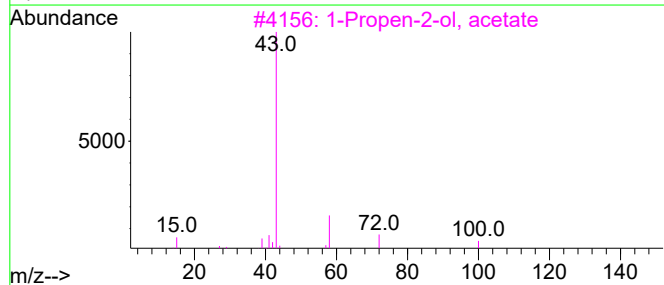
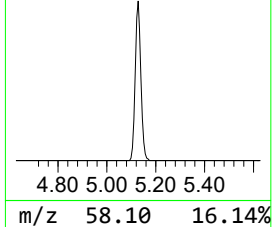
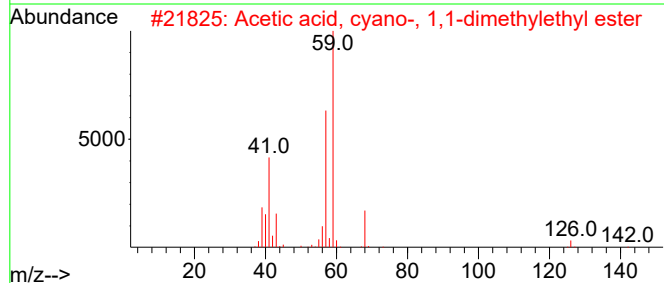
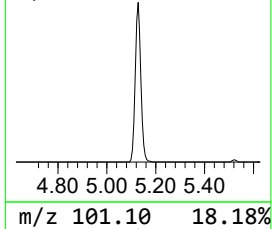
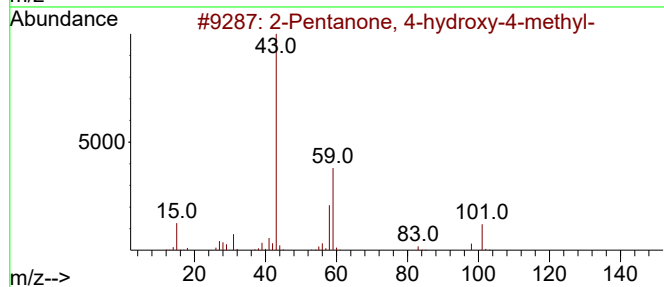
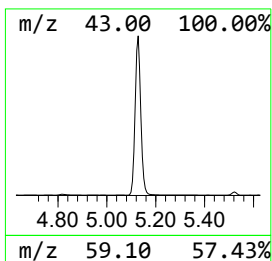
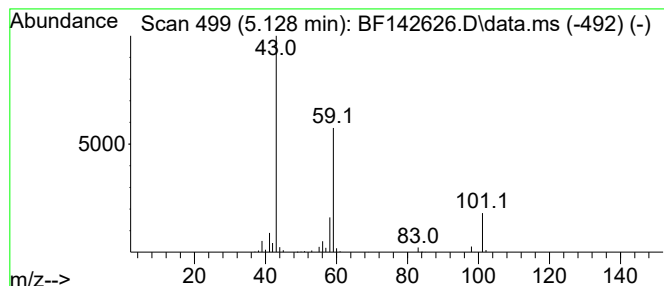
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD			R.T.
5.128	9.94 ng	338476	1,4-Dichlorobenzene-d4			6.892
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-		116	C6H12O2	000123-42-2	59
2	Acetic acid, cyano-, 1,1-dimethy...		141	C7H11NO2	001116-98-9	25
3	1-Propen-2-ol, acetate		100	C5H8O2	000108-22-5	12
4	Morpholine, 4-methyl-		101	C5H11NO	000109-02-4	9
5	2,3-Butanedione, monooxime		101	C4H7NO2	000057-71-6	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-...	5.128	9.9	ng	338476	1	6.892	681009	20.0