

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF082222\
 Data File : BF129937.D
 Acq On : 22 Aug 2022 12:52
 Operator : CG\JU
 Sample : PB146994BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB146994BL

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-BF081422.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BF129937.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	4.387	354	357	371	rBV	85939	129529	2.25%	0.392%
2	5.010	460	463	484	rBV	608876	950430	16.50%	2.879%
3	5.422	529	533	555	rBV	3957585	4031724	70.00%	12.211%
4	6.434	700	705	708	rBV	3977894	3978861	69.09%	12.051%
5	6.775	759	763	774	rVB	994053	828162	14.38%	2.508%
6	7.345	855	860	863	rBV	2561378	2644085	45.91%	8.008%
7	8.057	976	981	988	rBV	1248030	1132160	19.66%	3.429%
8	9.134	1158	1164	1167	rBV	5423001	5060946	87.88%	15.328%
9	9.810	1274	1279	1283	rBV	1581789	1340362	23.27%	4.060%
10	10.610	1409	1415	1418	rBV	3839186	3356878	58.29%	10.167%
11	11.298	1529	1532	1536	rBV	1572166	1378081	23.93%	4.174%
12	12.892	1797	1803	1806	rBV	6198840	5759235	100.00%	17.443%
13	13.945	1978	1982	1993	rBV	1376999	1288064	22.37%	3.901%
14	15.404	2222	2230	2237	rBV	895322	1138301	19.76%	3.448%

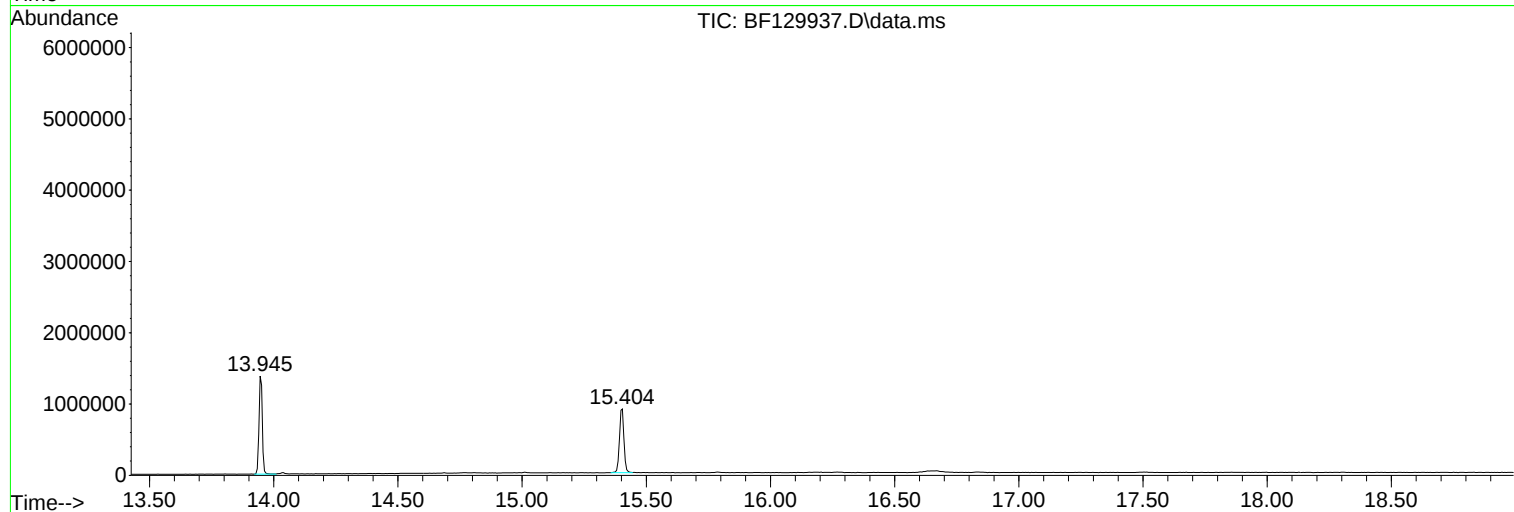
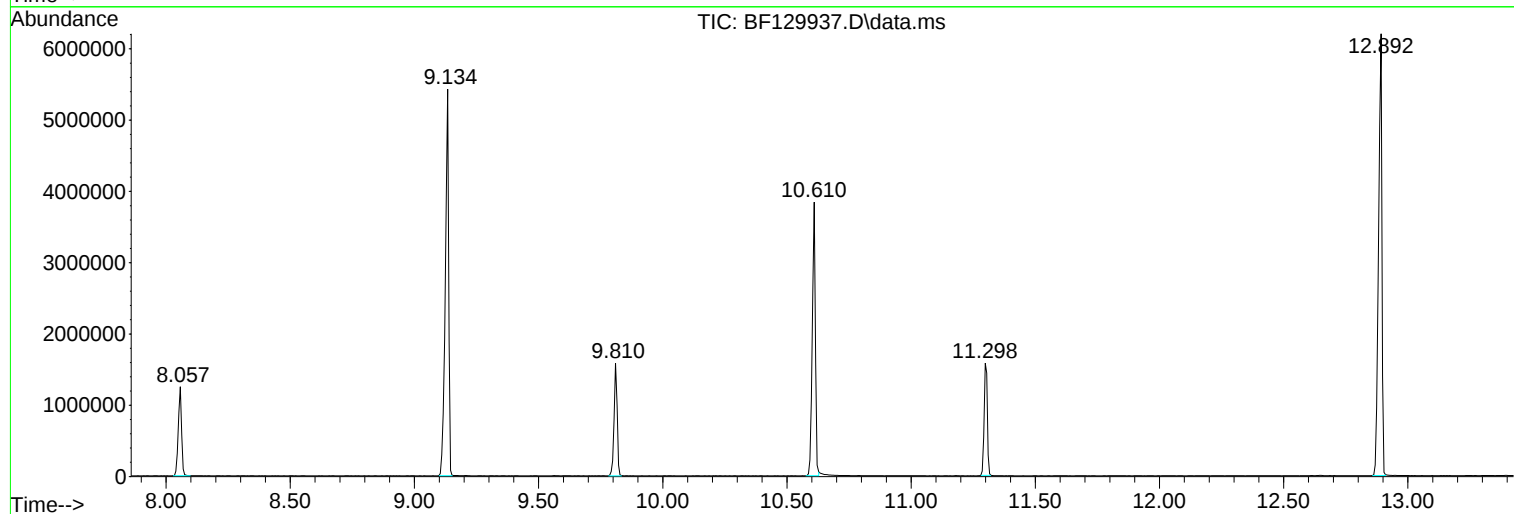
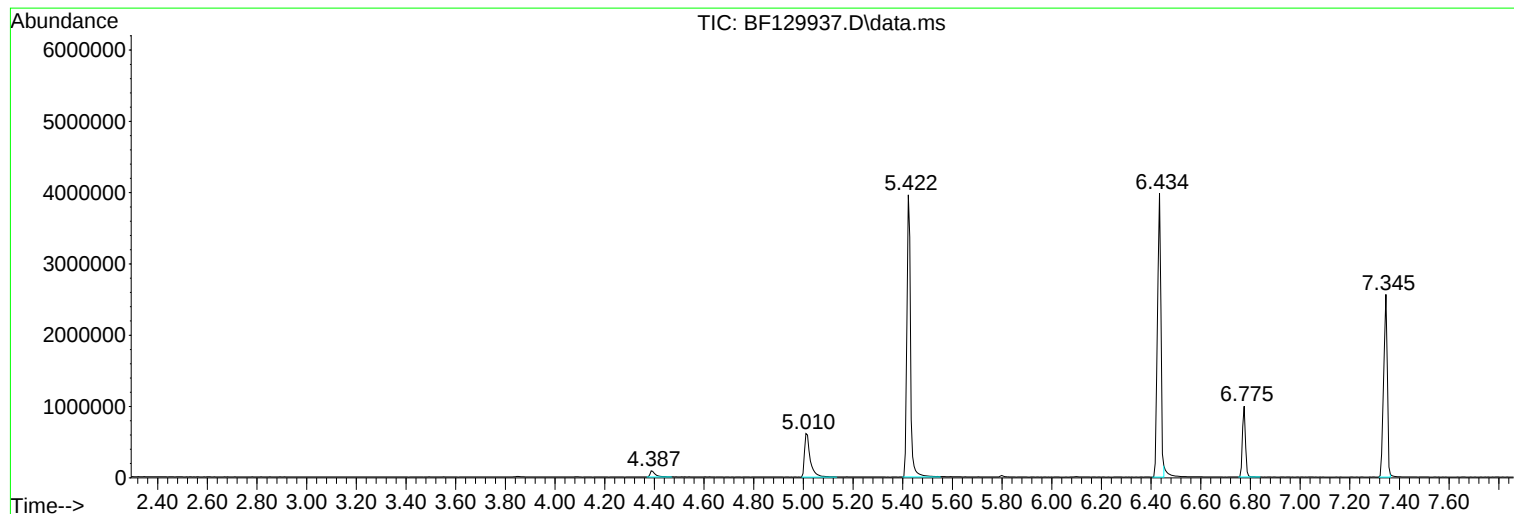
Sum of corrected areas: 33016818

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF081422.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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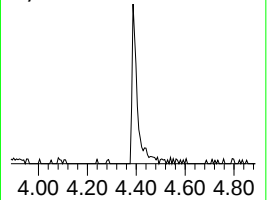
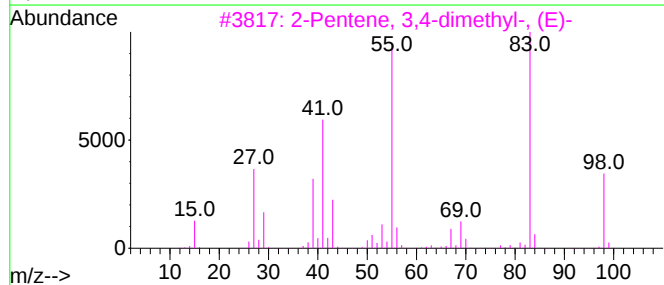
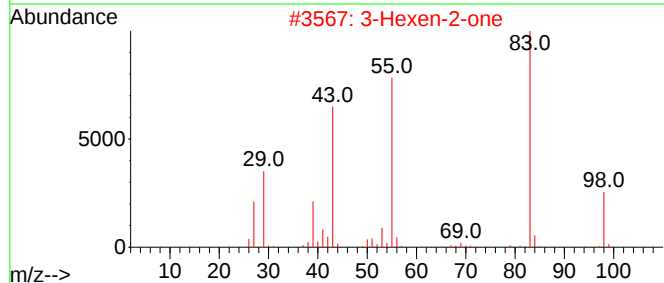
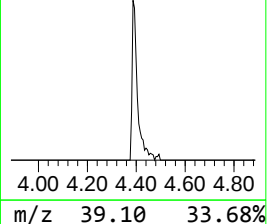
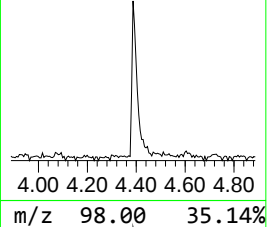
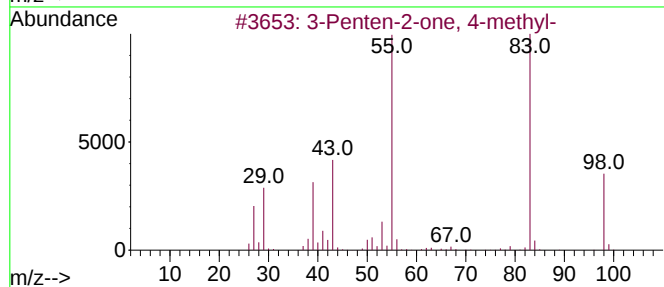
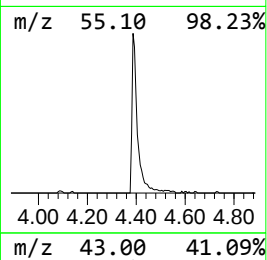
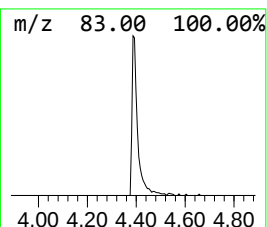
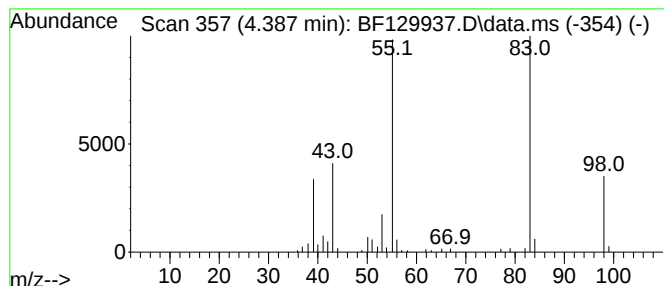
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF081422.M
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 3-Penten-2-one, 4-methyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.387	3.13 ng	129529	1,4-Dichlorobenzene-d4	6.775

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		3-Penten-2-one, 4-methyl-	98	C6H10O	000141-79-7	94
2		3-Hexen-2-one	98	C6H10O	000763-93-9	91
3		2-Pentene, 3,4-dimethyl-, (E)-	98	C7H14	004914-92-5	90
4		2-Pentene, 2,4-dimethyl-	98	C7H14	000625-65-0	86
5		2-Pentene, 4,4-dimethyl-, (E)-	98	C7H14	000690-08-4	86



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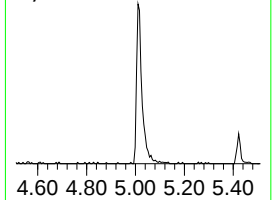
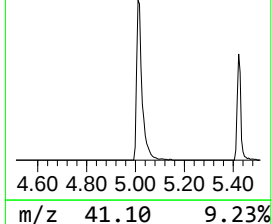
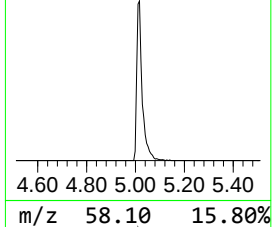
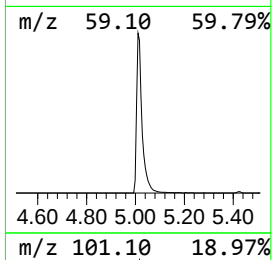
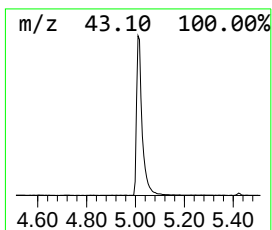
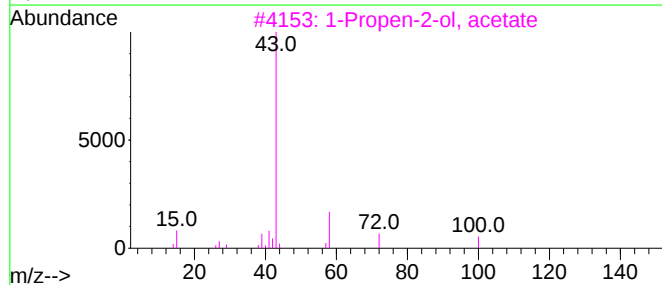
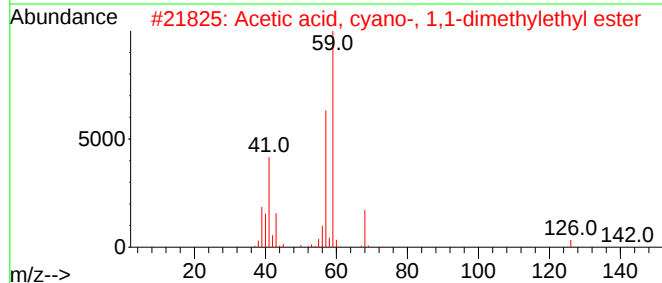
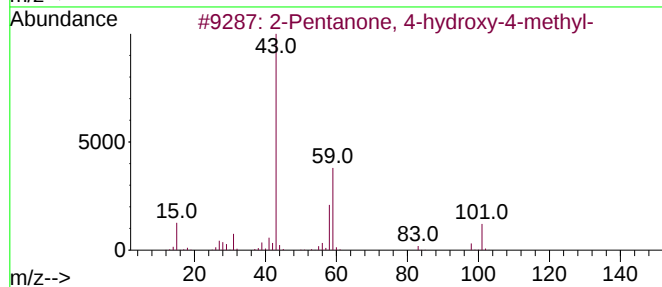
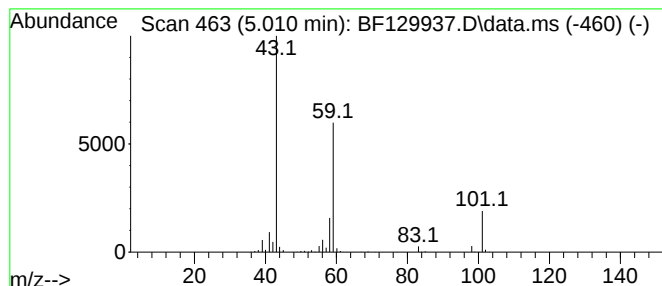
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Peak Number 2 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.010	22.95 ng	950430	1,4-Dichlorobenzene-d4	6.775

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	64
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
3-Penten-2-one,...	4.387	3.1	ng	129529	1	6.775	828162	20.0
2-Pentanone, 4-...	5.010	22.9	ng	950430	1	6.775	828162	20.0