

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF092822\
 Data File : BF130442.D
 Acq On : 28 Sep 2022 11:38
 Operator : CG\JU
 Sample : PB147922BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB147922BL

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-BF091422.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BF130442.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	4.175	351	355	370	rBB	68342	87446	2.22%	0.366%
2	4.840	463	468	483	rBV	417417	647676	16.46%	2.713%
3	5.257	534	539	542	rBV	2820698	3026080	76.89%	12.678%
4	6.287	709	714	717	rBV	2606674	3026346	76.90%	12.679%
5	6.640	771	774	778	rBV	778339	693913	17.63%	2.907%
6	7.210	865	871	874	rBV	1796408	2006834	50.99%	8.408%
7	7.928	987	993	996	rBV	938366	933528	23.72%	3.911%
8	9.004	1170	1176	1179	rBV	4211847	3935406	100.00%	16.487%
9	9.675	1284	1290	1294	rBV	1212660	1052711	26.75%	4.410%
10	10.469	1419	1425	1428	rBV	2671185	2503752	63.62%	10.489%
11	11.157	1538	1542	1545	rBV	1248135	1000879	25.43%	4.193%
12	12.745	1807	1812	1815	rBV	4305956	3703782	94.11%	15.517%
13	13.786	1985	1989	1997	rVB	818262	727344	18.48%	3.047%
14	15.174	2220	2225	2230	rBV	473446	523760	13.31%	2.194%

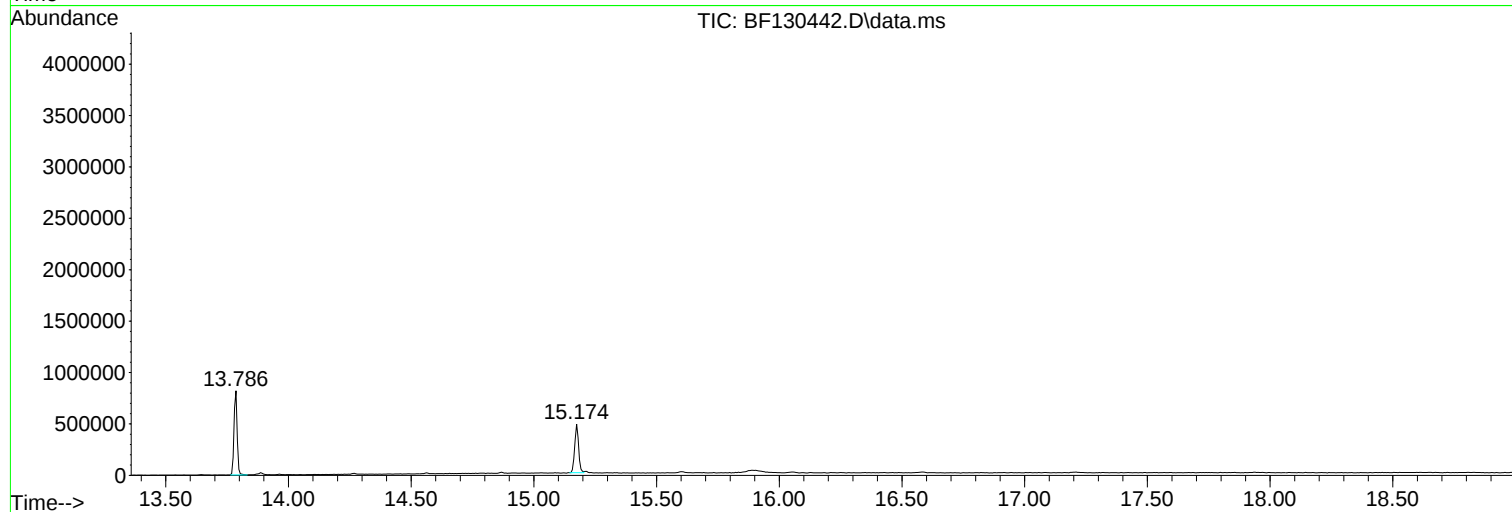
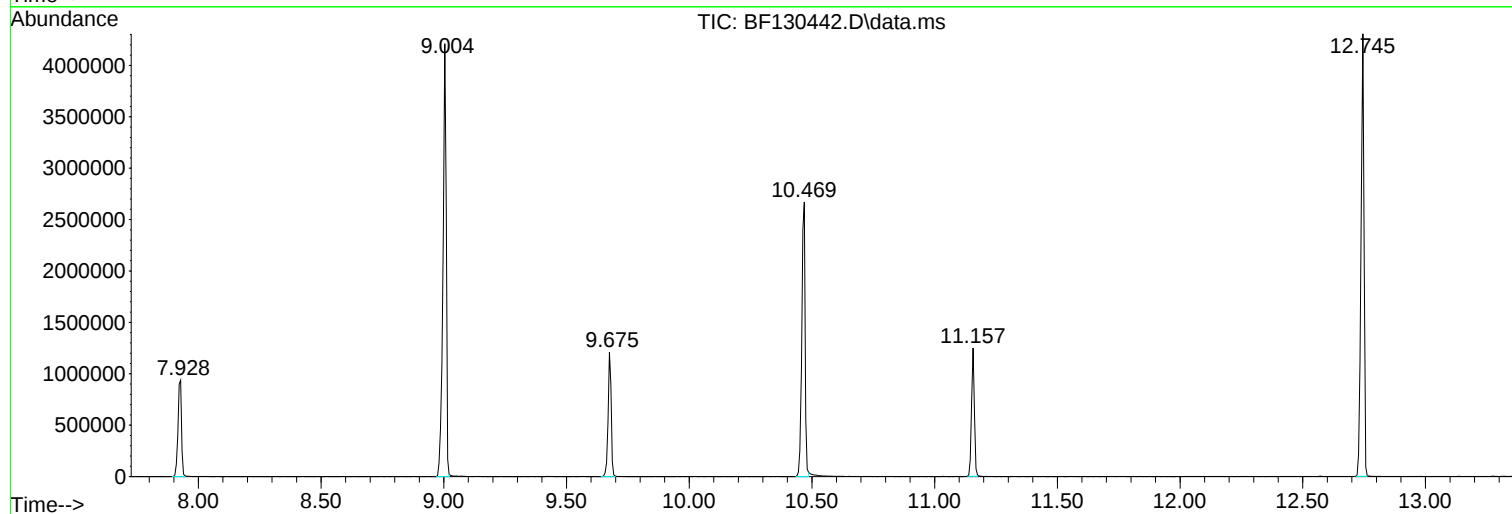
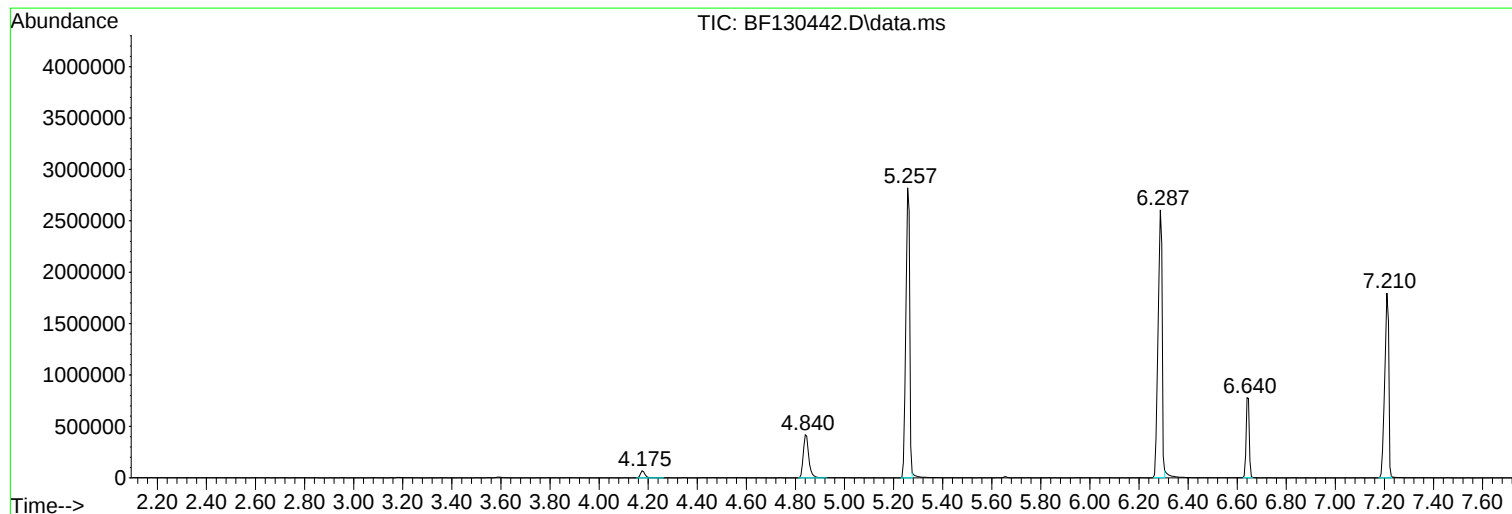
Sum of corrected areas: 23869457

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF091422.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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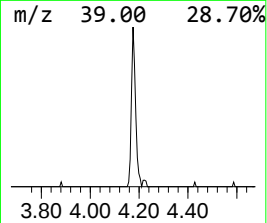
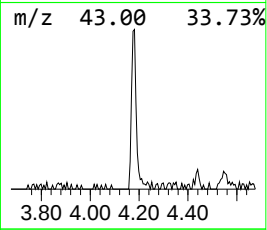
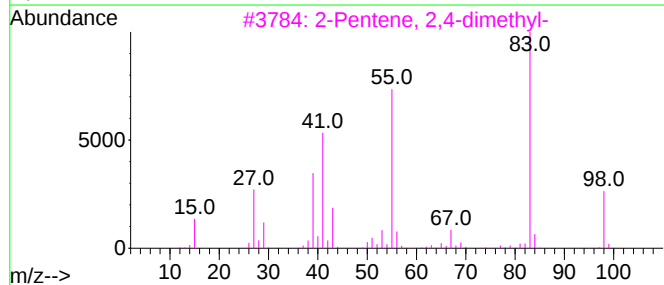
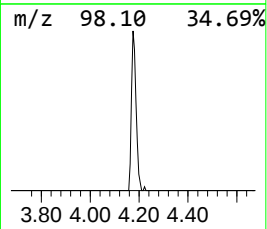
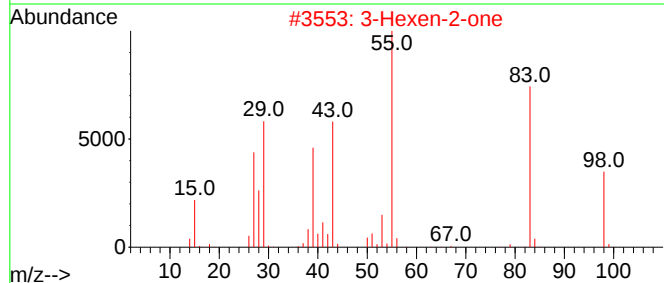
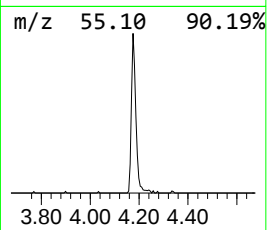
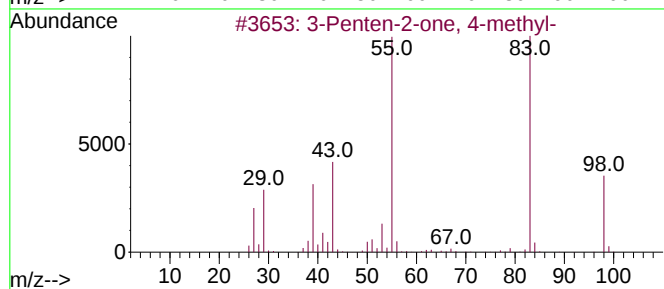
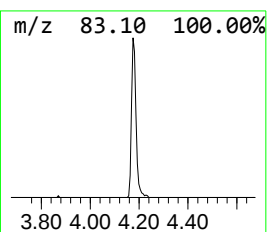
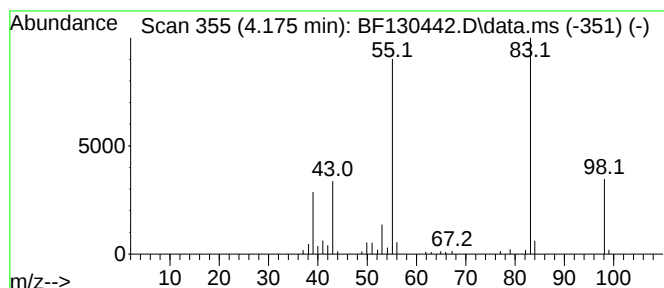
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TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.175	2.52 ng	87446	1,4-Dichlorobenzene-d4	6.646

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, 2,4-dimethyl-	98	C7H14	000625-65-0	87
4		2-Pentene, 3,4-dimethyl-, (E)-	98	C7H14	004914-92-5	86
5		2-Pentene, 4,4-dimethyl-, (E)-	98	C7H14	000690-08-4	86



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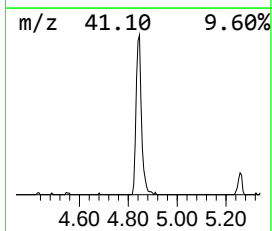
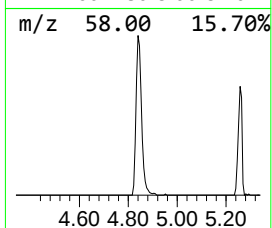
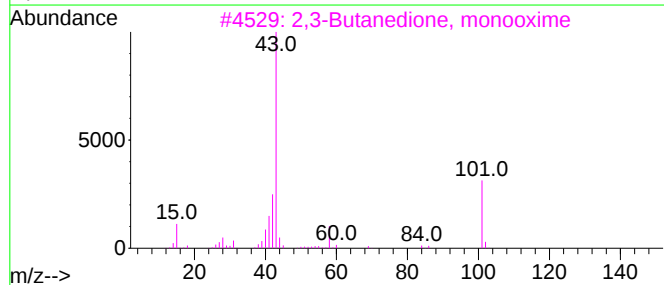
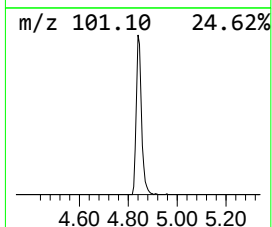
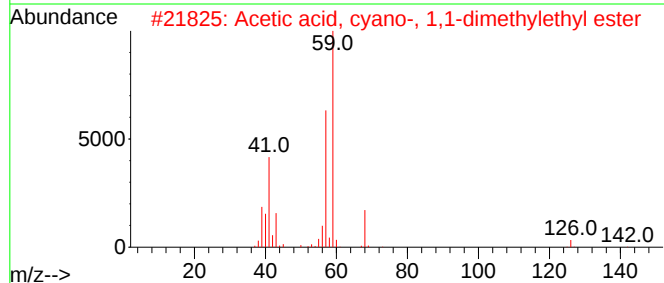
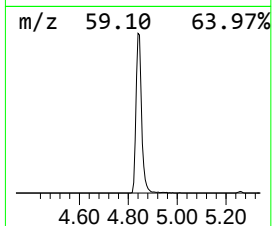
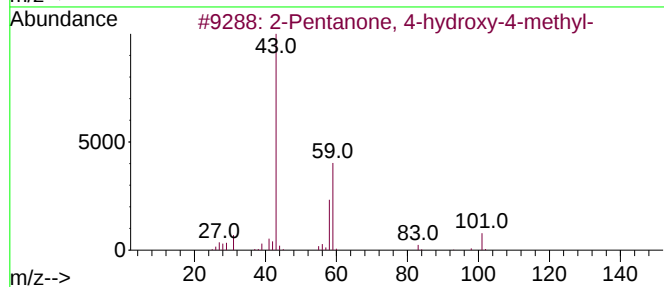
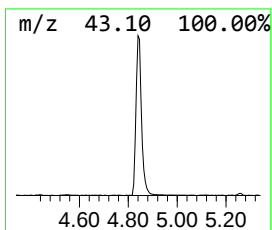
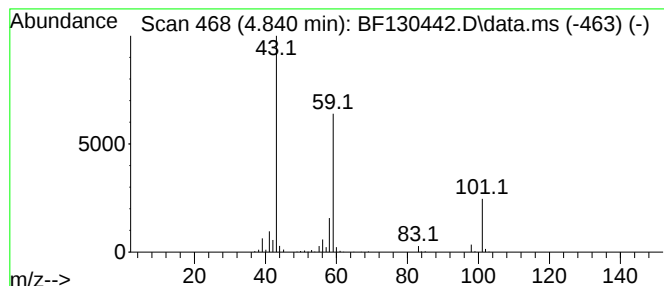
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Peak Number 2 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.840	18.67 ng	647676	1,4-Dichlorobenzene-d4	6.646

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	25
3		2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	17
4		Acetone	58	C3H6O	000067-64-1	10
5		1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	10



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
3-Penten-2-one,...	4.175	2.5	ng	87446	1	6.646	693913	20.0
2-Pentanone, 4-...	4.840	18.7	ng	647676	1	6.646	693913	20.0