

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF061923\
 Data File : BF133831.D
 Acq On : 19 Jun 2023 13:35
 Operator : CG\JU
 Sample : PB153132BL
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB153132BL

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-BF060923.M
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BF133831.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	4.504	407	411	418	rBV	92502	108723	2.52%	0.392%
2	5.110	509	514	524	rBV	748100	1044161	24.21%	3.761%
3	5.504	575	581	584	rBV	3474083	3480038	80.70%	12.534%
4	6.498	744	750	753	rBV	3149883	3552307	82.38%	12.794%
5	6.857	807	811	815	rBV	978621	752793	17.46%	2.711%
6	7.422	901	907	910	rBV	2442055	2408667	55.86%	8.675%
7	8.139	1025	1029	1032	rBV	1092084	984273	22.82%	3.545%
8	9.216	1208	1212	1215	rBV	4785435	4267561	98.96%	15.371%
9	9.898	1324	1328	1331	rBV	1142427	1058185	24.54%	3.811%
10	10.692	1458	1463	1466	rBV	2954659	2962362	68.70%	10.670%
11	11.386	1577	1581	1585	rBV	1447873	1301322	30.18%	4.687%
12	12.992	1848	1854	1857	rBV	4524366	4312279	100.00%	15.532%
13	14.045	2028	2033	2041	rBV	979143	873485	20.26%	3.146%
14	15.551	2284	2289	2295	rBV	491186	658282	15.27%	2.371%

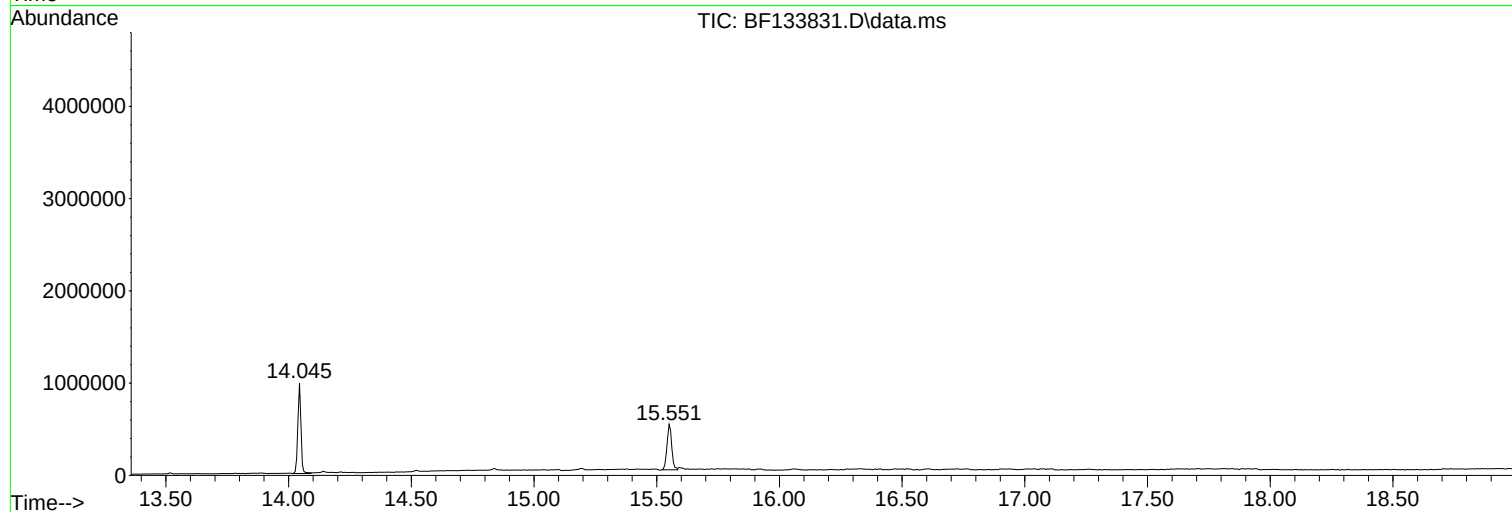
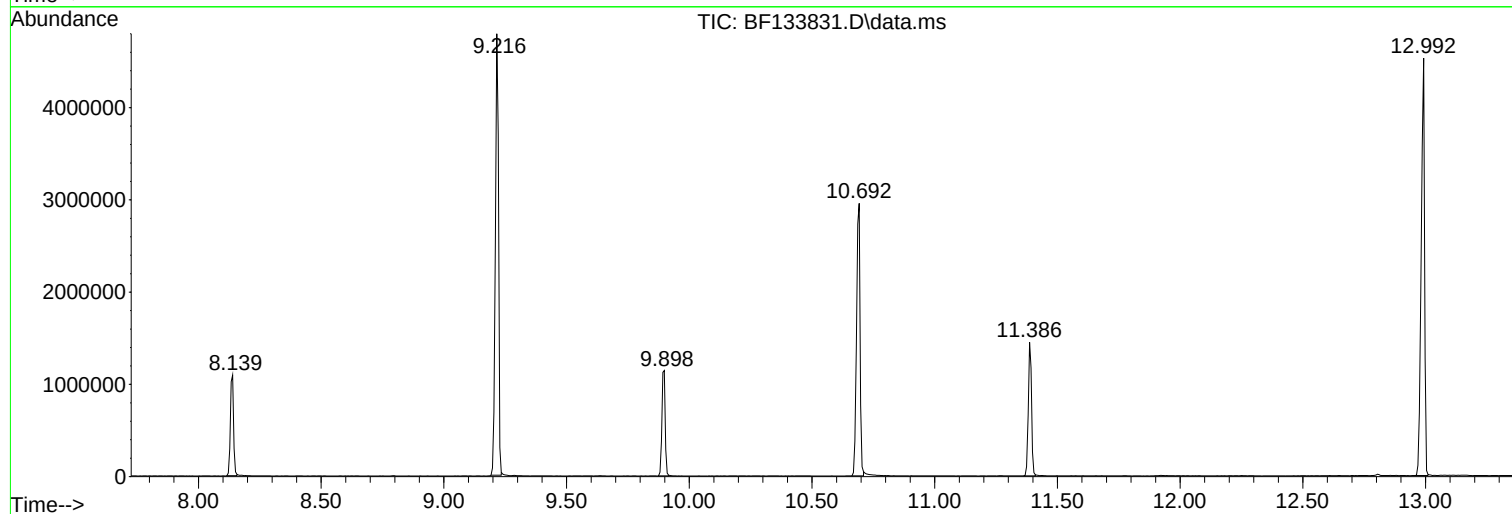
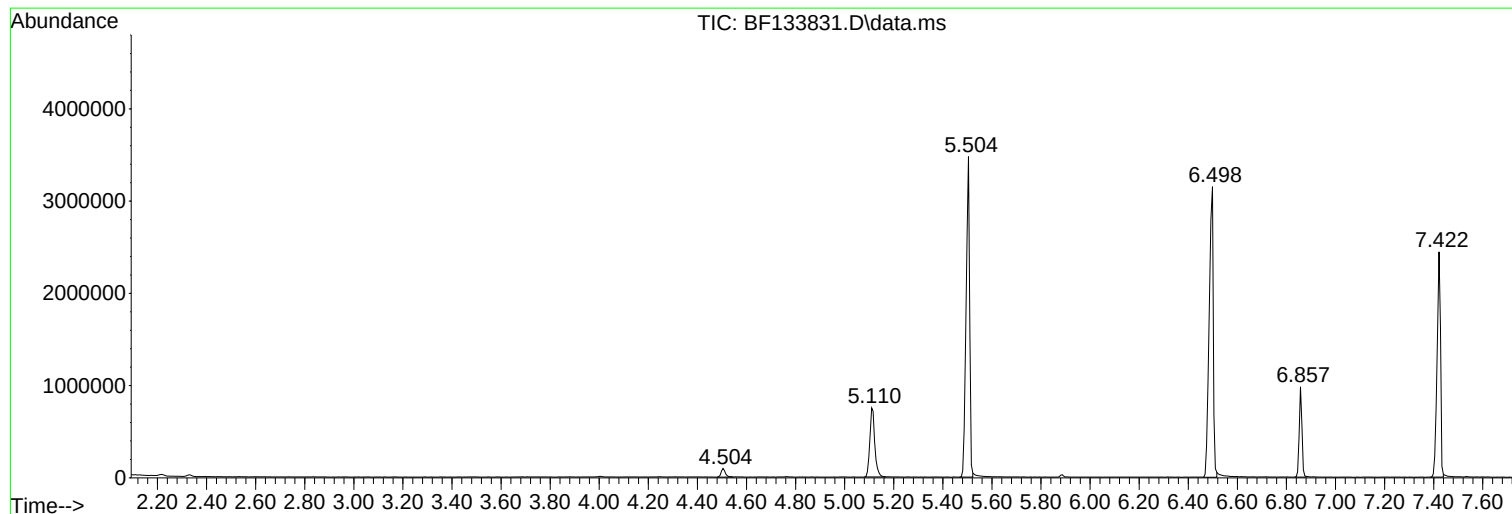
Sum of corrected areas: 27764438

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF060923.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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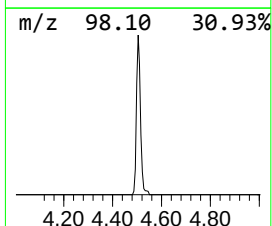
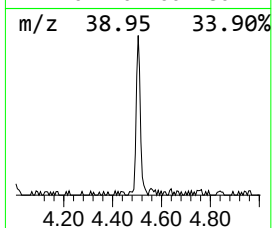
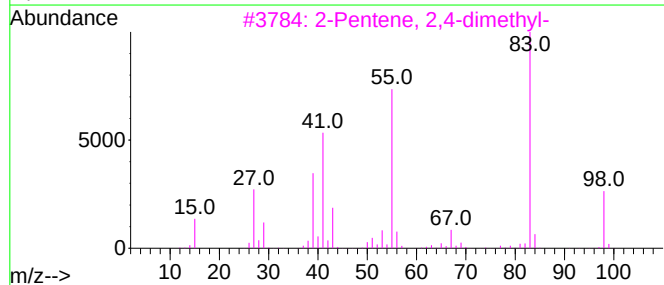
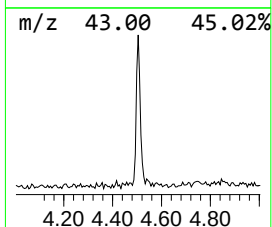
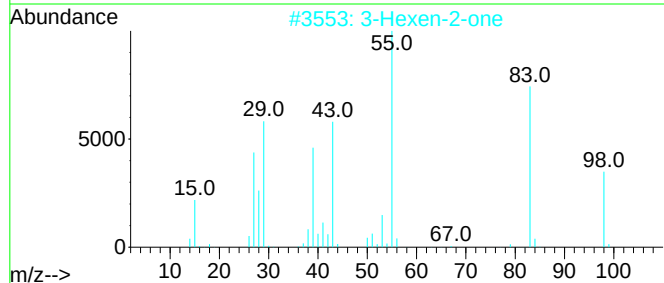
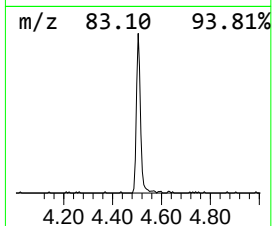
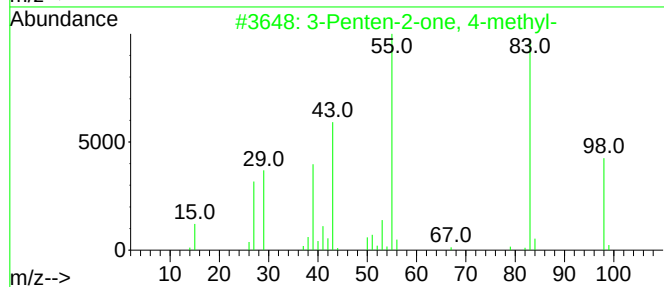
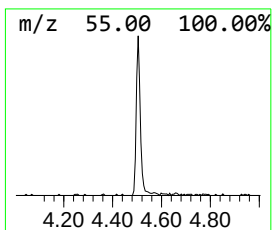
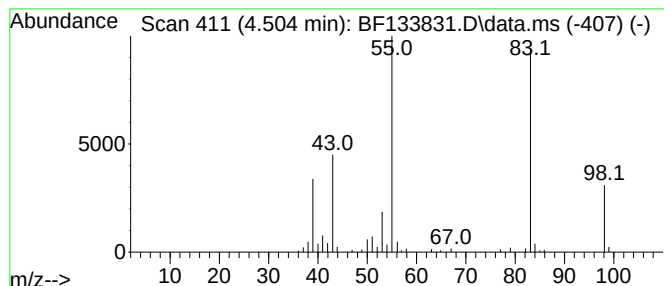
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF060923.M
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.504	2.89 ng	108723	1,4-Dichlorobenzene-d4	6.857

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		3-Penten-2-one, 4-methyl-	98	C6H10O	000141-79-7	91
2		3-Hexen-2-one	98	C6H10O	000763-93-9	91
3		2-Pentene, 2,4-dimethyl-	98	C7H14	000625-65-0	80
4		2-Pentene, 3,4-dimethyl-	98	C7H14	024910-63-2	80
5		Cyclohexane, methyl-	98	C7H14	000108-87-2	72



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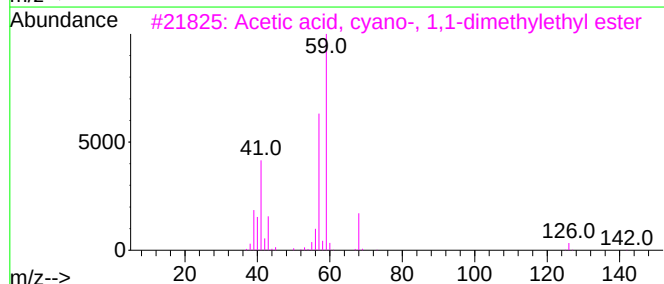
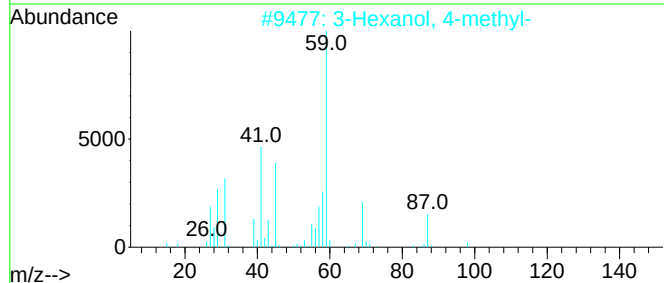
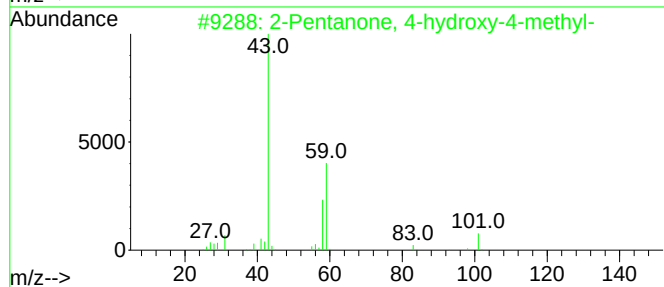
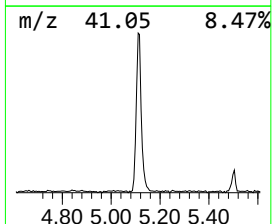
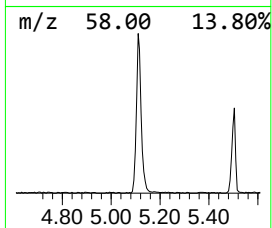
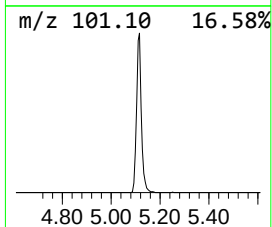
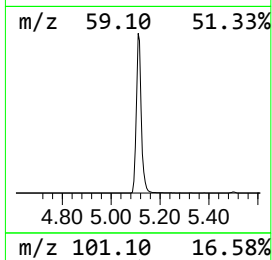
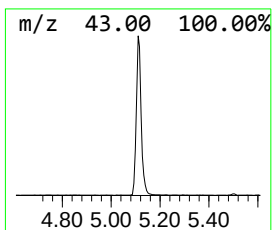
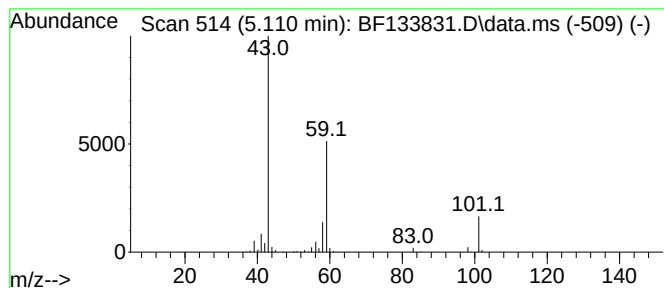
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Peak Number 2 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.110	27.74 ng	1044160	1,4-Dichlorobenzene-d4	6.857

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50		
2	3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	33		
3	Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	25		
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
3-Penten-2-one,...	4.504	2.9	ng	108723	1	6.857	752793	20.0
2-Pentanone, 4-...	5.110	27.7	ng	1044160	1	6.857	752793	20.0