

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111722\
 Data File : BP012663.D
 Acq On : 17 Nov 2022 16:07
 Operator : CG/JU
 Sample : PB148901BL
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 PB148901BL

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_P\Methods\8270E-BP111722.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BP012663.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	4.505	253	258	266	rBV	258548	364512	1.46%	0.360%
2	5.116	357	362	373	rVB	1398185	1998889	8.02%	1.974%
3	5.587	435	442	455	rBV	5456917	8057975	32.32%	7.959%
4	7.193	707	715	726	rBV	5504083	8647753	34.69%	8.541%
5	8.046	853	860	867	rVB	1006538	1578345	6.33%	1.559%
6	9.210	1050	1058	1068	rBV	3050742	4977463	19.96%	4.916%
7	10.869	1330	1340	1348	rBV	1473543	2469482	9.91%	2.439%
8	13.310	1748	1755	1764	rBV	9368856	14157549	56.79%	13.983%
9	14.681	1981	1988	1996	rBV	2554396	3753044	15.05%	3.707%
10	16.169	2234	2241	2264	rBV	8591219	12154090	48.75%	12.004%
11	17.439	2450	2457	2469	rVB	3412331	4892224	19.62%	4.832%
12	19.986	2884	2890	2895	rBV	20132982	24931416	100.00%	24.624%
13	21.521	3145	3151	3168	rVB	4657970	6292520	25.24%	6.215%
14	24.051	3574	3581	3592	rVB2	3407888	6972039	27.96%	6.886%

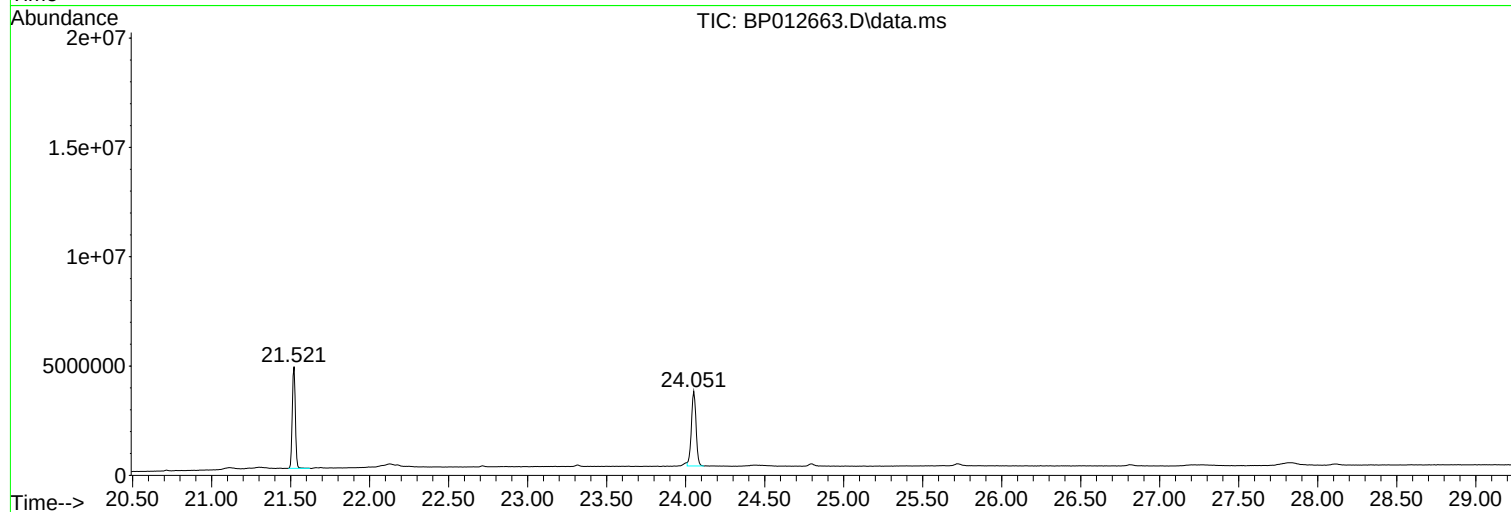
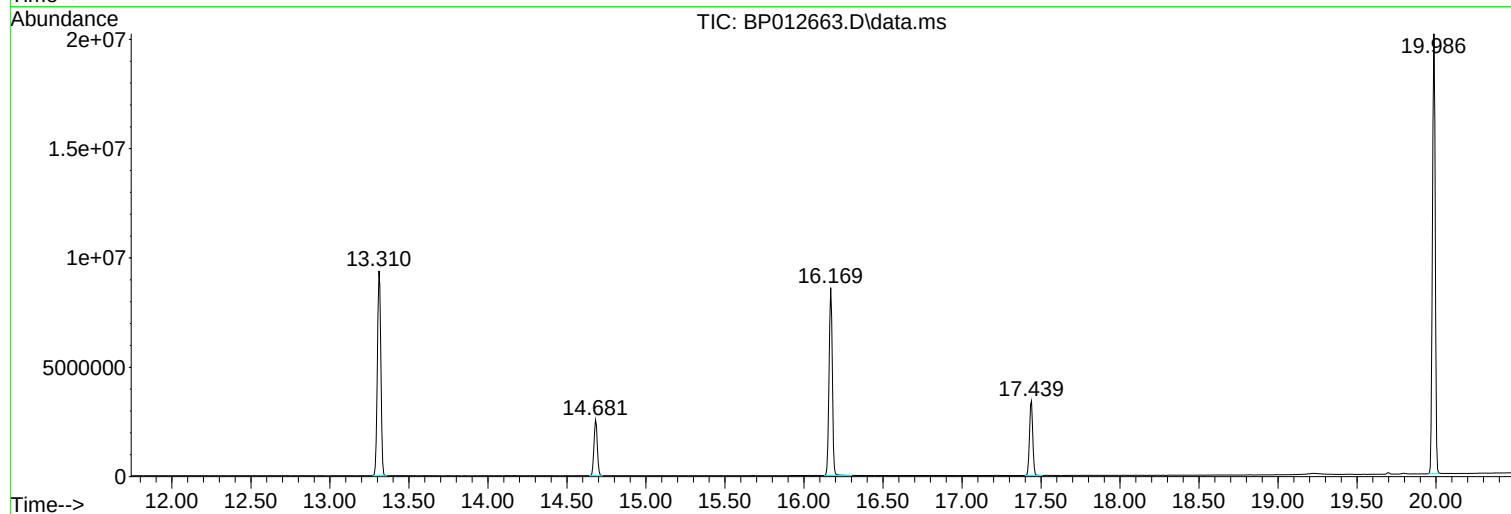
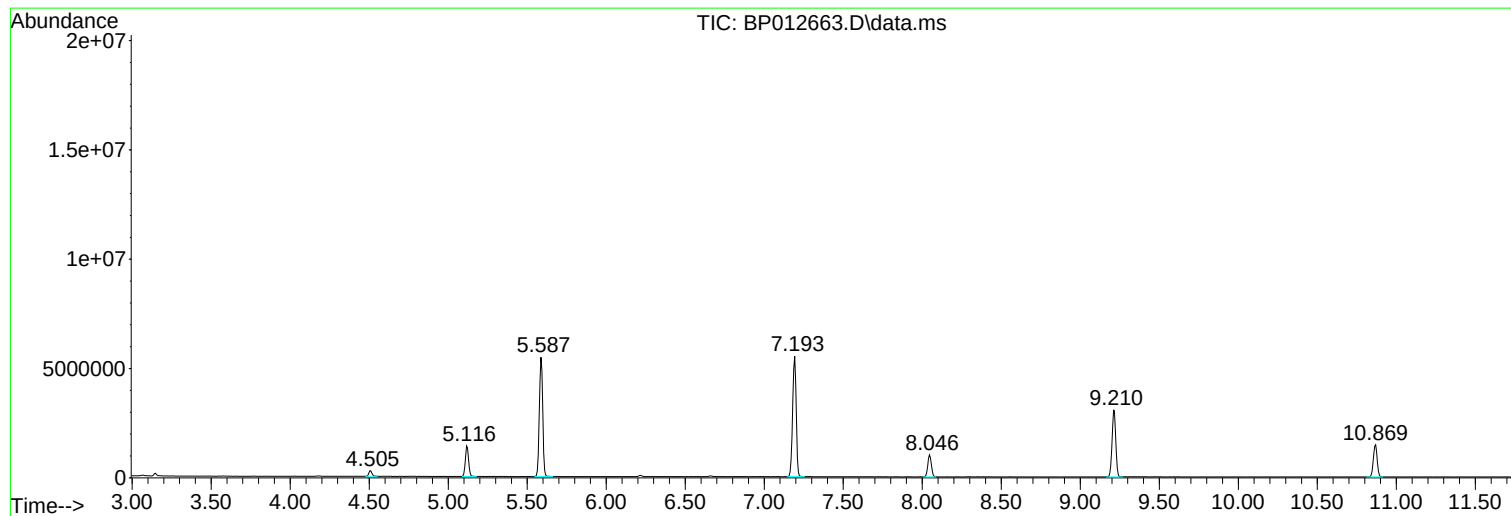
Sum of corrected areas: 101247301

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP111722.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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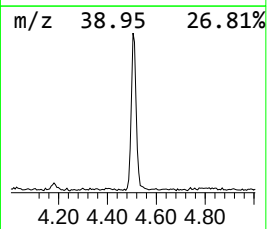
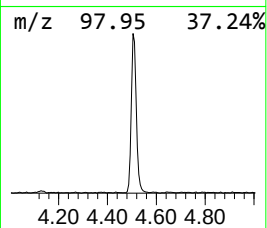
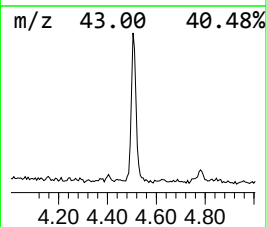
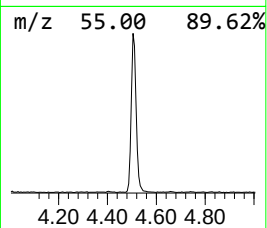
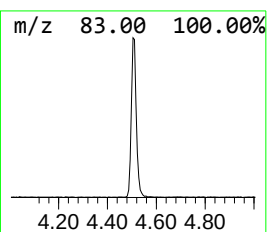
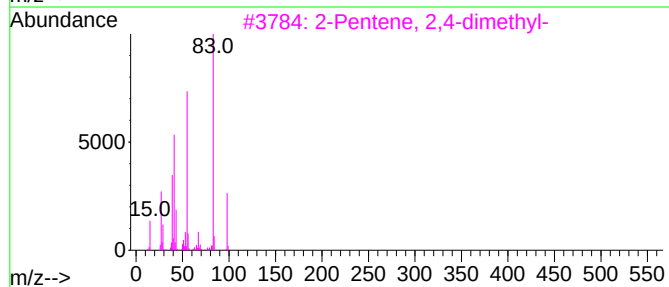
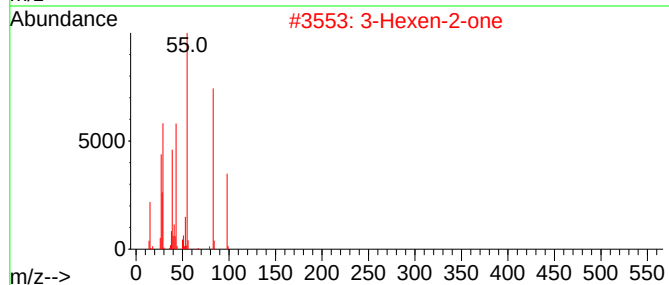
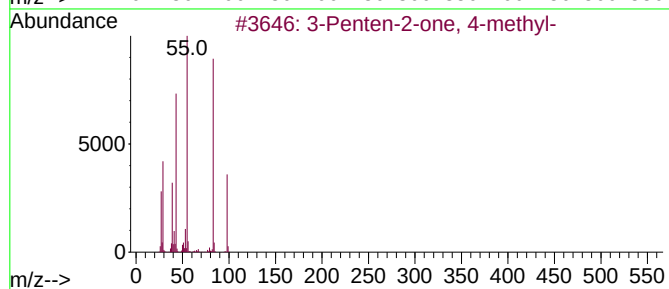
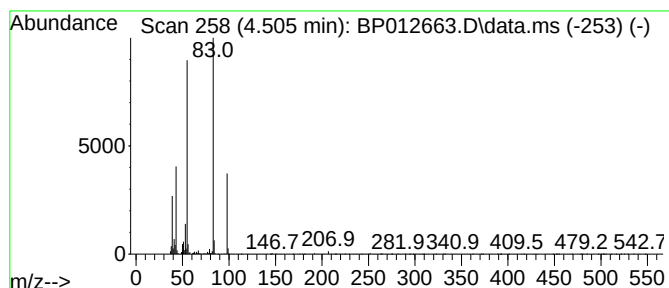
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP111722.M
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.505	4.62 ng	364512	1,4-Dichlorobenzene-d4	8.046

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3-Penten-2-one, 4-methyl-	98	C6H10O	000141-79-7	90
2			3-Hexen-2-one	98	C6H10O	000763-93-9	90
3			2-Pentene, 2,4-dimethyl-	98	C7H14	000625-65-0	87
4			2-Pentene, 4,4-dimethyl-, (E)-	98	C7H14	000690-08-4	86
5			2-Pentene, 3,4-dimethyl-, (E)-	98	C7H14	004914-92-5	86



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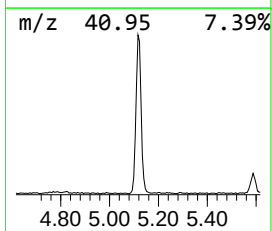
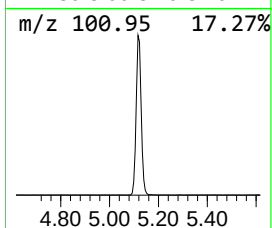
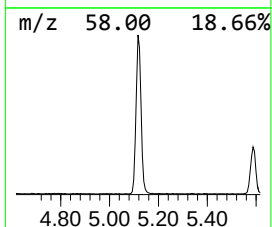
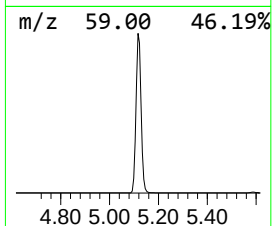
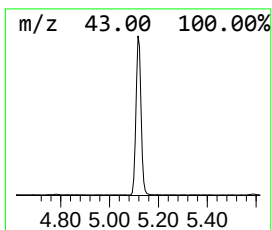
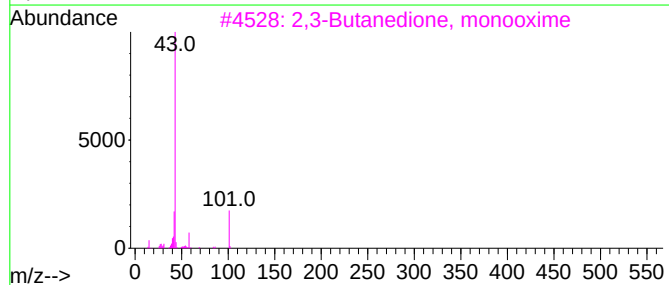
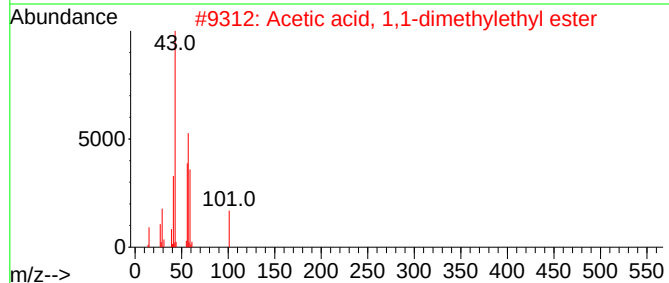
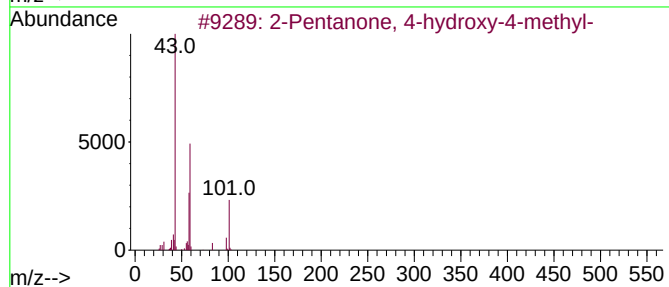
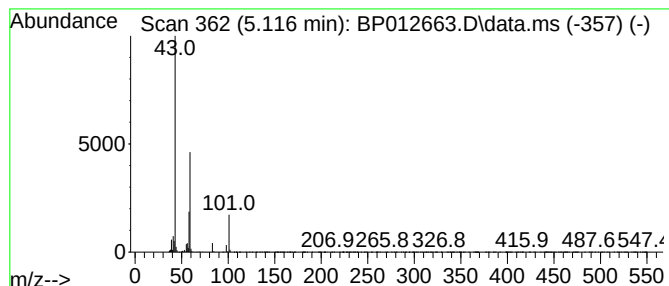
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Peak Number 2 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD			R.T.
5.116	25.33 ng	1998890	1,4-Dichlorobenzene-d4			8.046
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-		116	C6H12O2	000123-42-2	83
2	Acetic acid, 1,1-dimethylethyl e...		116	C6H12O2	000540-88-5	38
3	2,3-Butanedione, monooxime		101	C4H7NO2	000057-71-6	16
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,...	4.505	4.6	ng	364512	1	8.046	1578350	20.0
2-Pentanone, 4-...	5.116	25.3	ng	1998890	1	8.046	1578350	20.0