

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111722\
 Data File : BP012664.D
 Acq On : 17 Nov 2022 16:43
 Operator : CG/JU
 Sample : PB149128BL
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 PB149128BL

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: BP012664.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	307870	417695	2.00%	0.448%
2	5.117	356	362	372	rVB	1610595	2226508	10.66%	2.385%
3	5.587	435	442	454	rBV	5963366	8662792	41.47%	9.281%
4	7.187	706	714	727	rBV	5653771	9268145	44.37%	9.930%
5	8.040	853	859	868	rVB	1056852	1703973	8.16%	1.826%
6	9.210	1048	1058	1069	rBV	3278459	5274085	25.25%	5.651%
7	10.863	1329	1339	1352	rBV	1521726	2592671	12.41%	2.778%
8	13.304	1746	1754	1767	rBV	9318751	13943576	66.75%	14.939%
9	14.675	1980	1987	1995	rBV	2580268	3677871	17.61%	3.940%
10	16.163	2233	2240	2261	rVB	7391843	10273673	49.18%	11.007%
11	17.433	2449	2456	2465	rBV	3101742	4490249	21.50%	4.811%
12	19.980	2883	2889	2898	rBV	16779525	20888135	100.00%	22.380%
13	21.516	3144	3150	3161	rVB	3672650	4939231	23.65%	5.292%
14	24.039	3572	3579	3589	rVB2	2359099	4976705	23.83%	5.332%

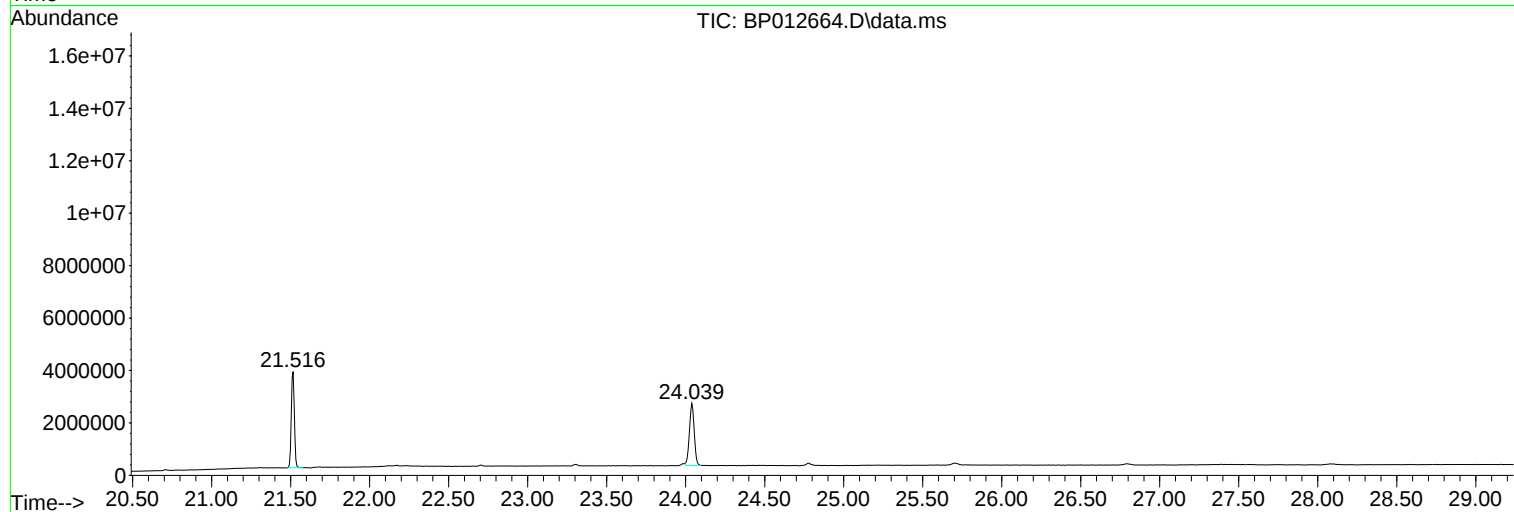
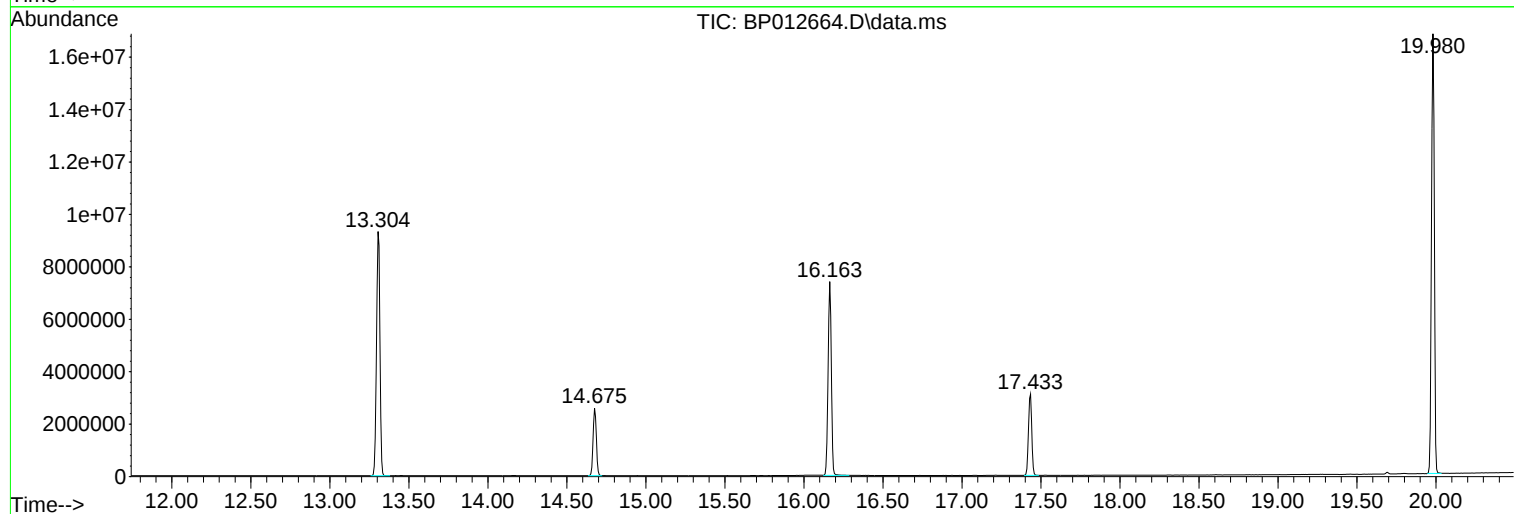
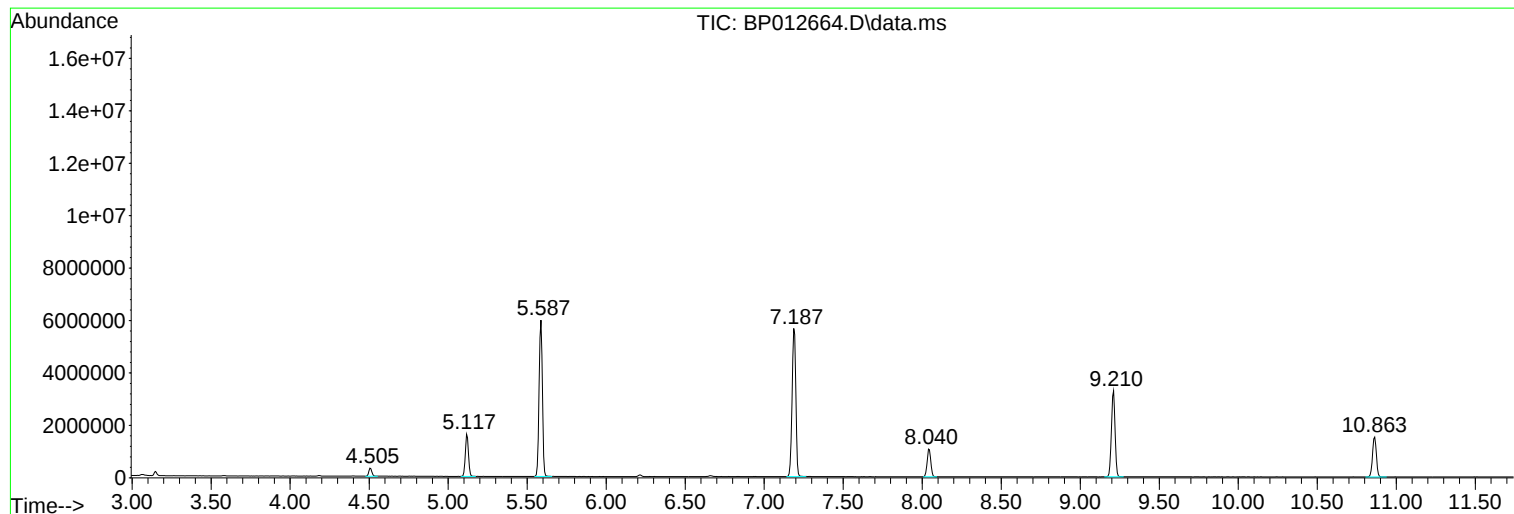
Sum of corrected areas: 93335309

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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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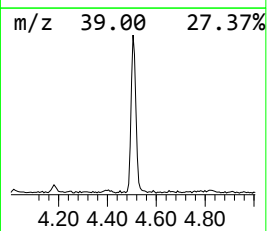
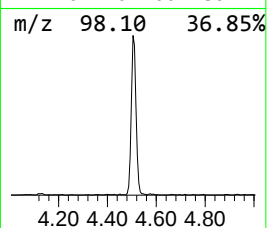
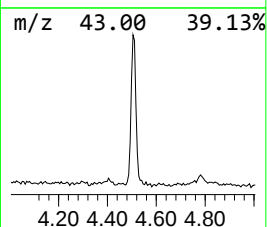
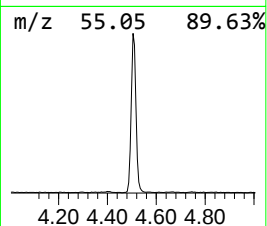
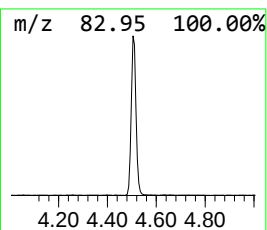
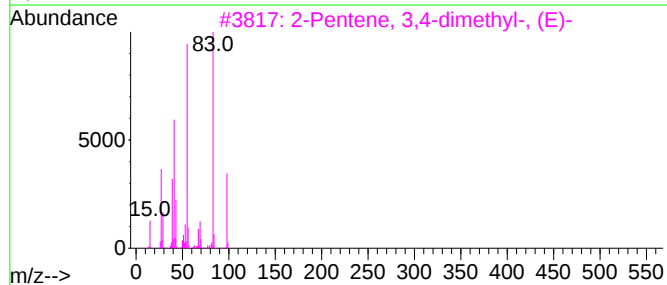
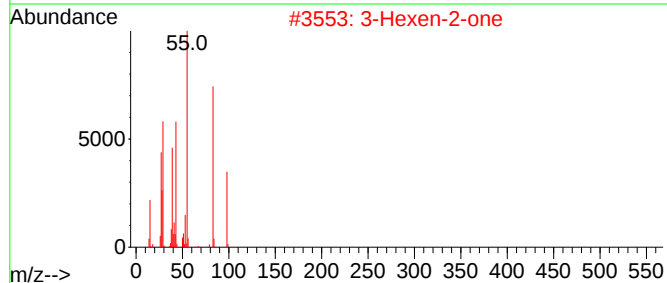
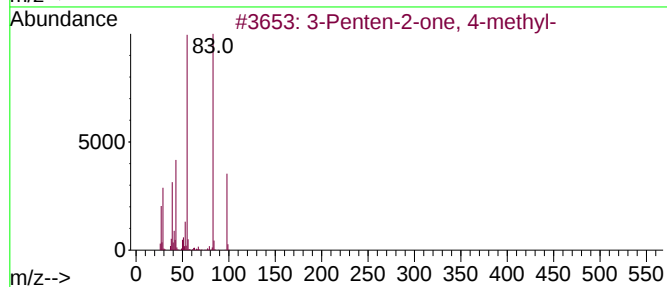
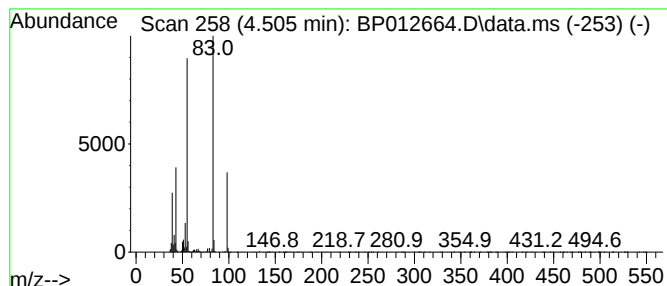
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP111722.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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	4.90 ng	417695	1,4-Dichlorobenzene-d4	8.040

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3-Penten-2-one, 4-methyl-	98	C6H10O	000141-79-7	95
2			3-Hexen-2-one	98	C6H10O	000763-93-9	91
3			2-Pentene, 3,4-dimethyl-, (E)-	98	C7H14	004914-92-5	86
4			2-Pentene, 2,4-dimethyl-	98	C7H14	000625-65-0	80
5			Furan, 2-methoxy-	98	C5H6O2	025414-22-6	78



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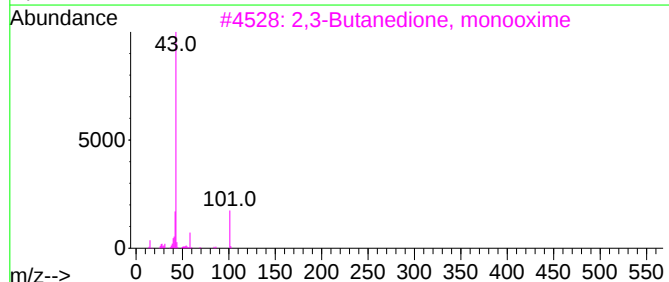
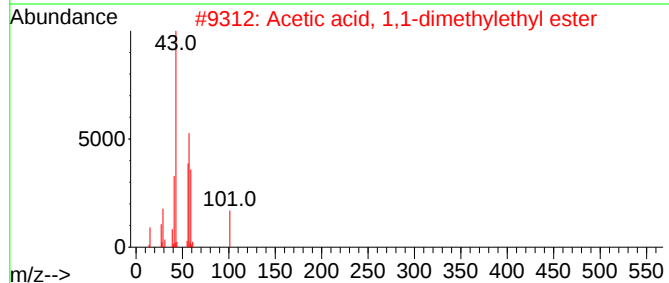
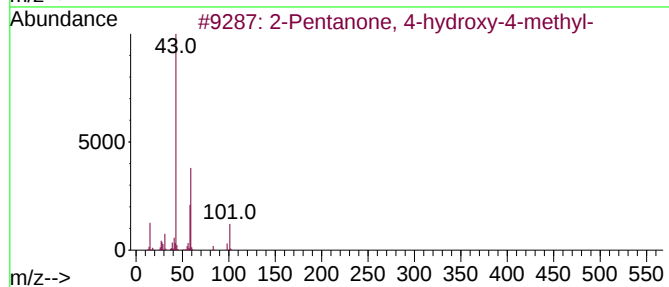
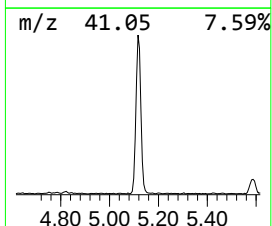
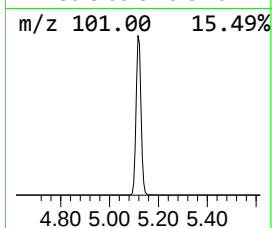
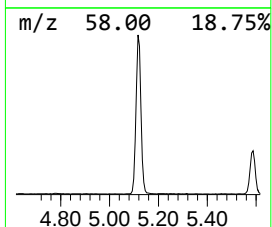
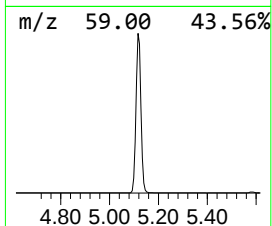
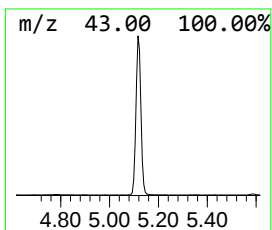
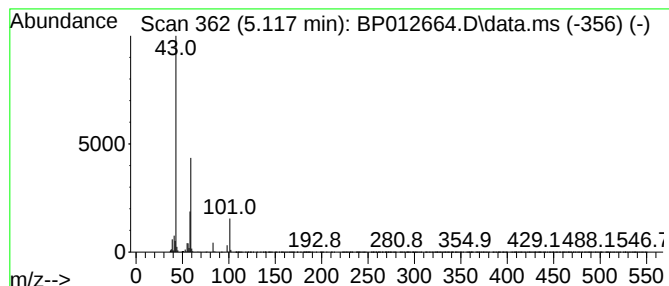
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TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.117	26.13 ng	2226510	1,4-Dichlorobenzene-d4	8.040

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	12
4			Acetone	58	C3H6O	000067-64-1	10
5			5-Hexen-2-one	98	C6H10O	000109-49-9	10



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
3-Penten-2-one,...	4.505	4.9	ng	417695	1	8.040	1703970	20.0
2-Pentanone, 4-...	5.117	26.1	ng	2226510	1	8.040	1703970	20.0