

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP030123\
 Data File : BP013914.D
 Acq On : 01 Mar 2023 11:27
 Operator : CG/JU
 Sample : PB151104BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 PB151104BL

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-BP021623.M
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BP013914.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	4.546	225	231	239	rBV	71360	96900	1.24%	0.287%
2	5.164	330	336	348	rBV	1112178	1563529	20.01%	4.633%
3	5.634	409	416	433	rBV	2129557	3066913	39.26%	9.087%
4	7.252	680	691	705	rBV	1905866	3064956	39.23%	9.081%
5	8.093	828	834	844	rBV	404667	639713	8.19%	1.895%
6	9.269	1026	1034	1051	rBV	1183814	1958053	25.06%	5.802%
7	10.922	1303	1315	1326	rBV	489980	832909	10.66%	2.468%
8	13.357	1719	1729	1743	rBV	3050739	4552640	58.28%	13.489%
9	14.734	1956	1963	1975	rBV	704656	1087158	13.92%	3.221%
10	16.222	2209	2216	2237	rBV	2622283	3887394	49.76%	11.518%
11	17.487	2424	2431	2441	rBV	973256	1396045	17.87%	4.136%
12	20.022	2856	2862	2873	rBV	6533788	7812101	100.00%	23.147%
13	21.563	3119	3124	3130	rBV	1293689	1765792	22.60%	5.232%
14	24.110	3550	3557	3567	rVB2	917131	2025522	25.93%	6.002%

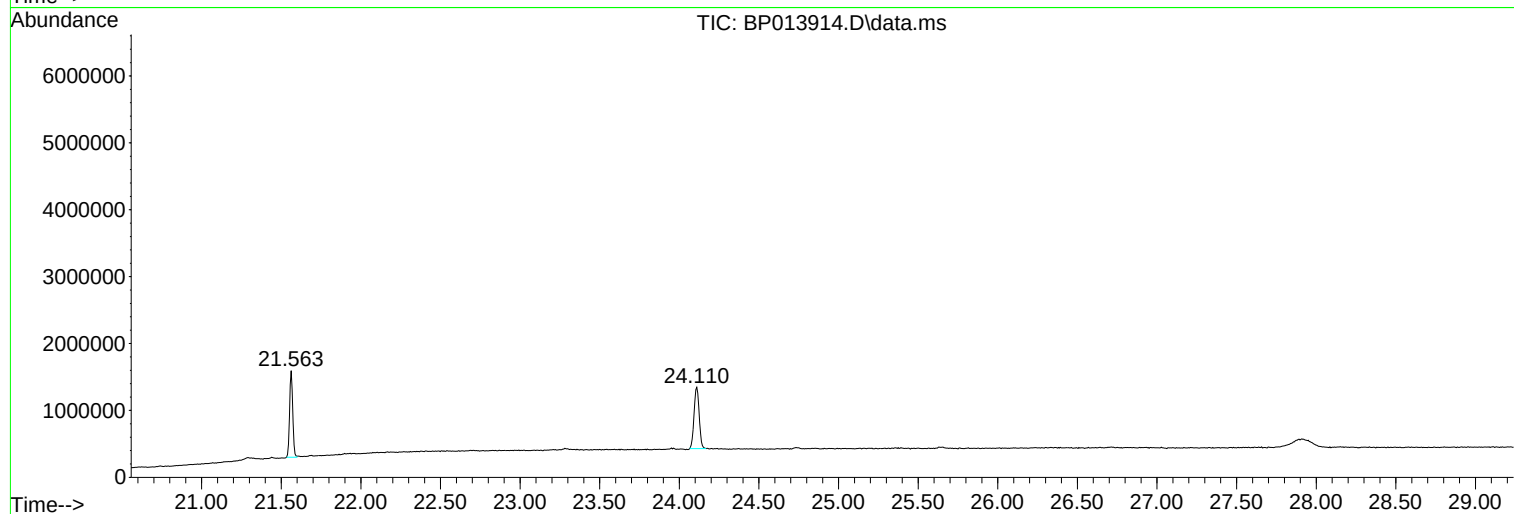
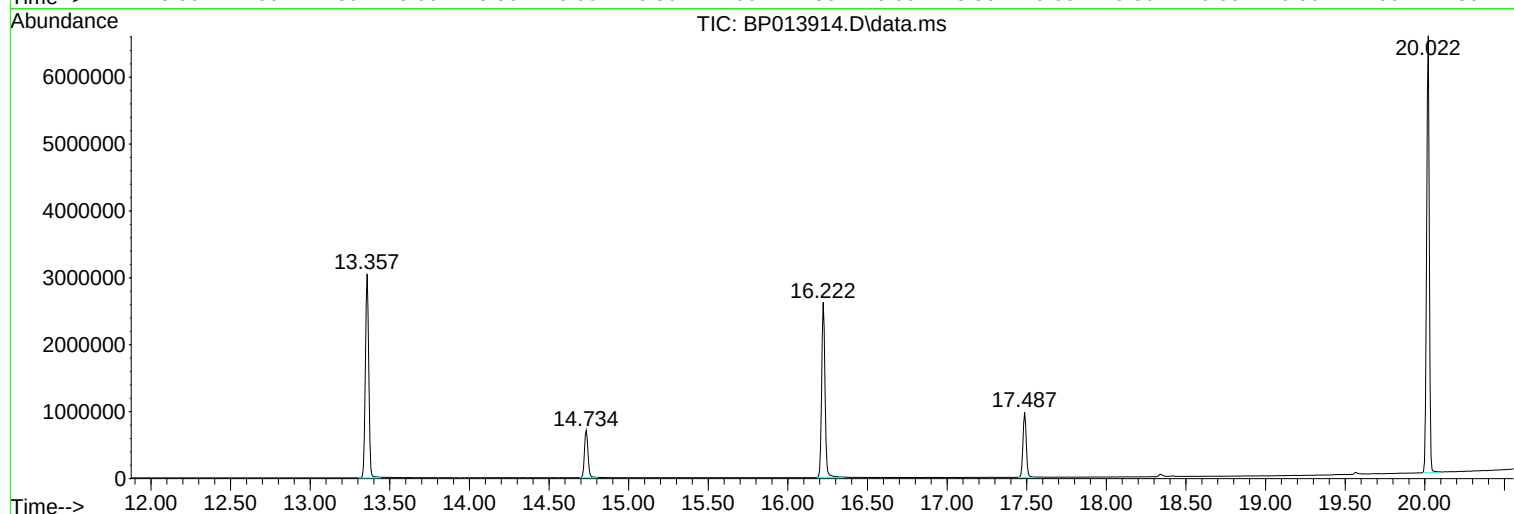
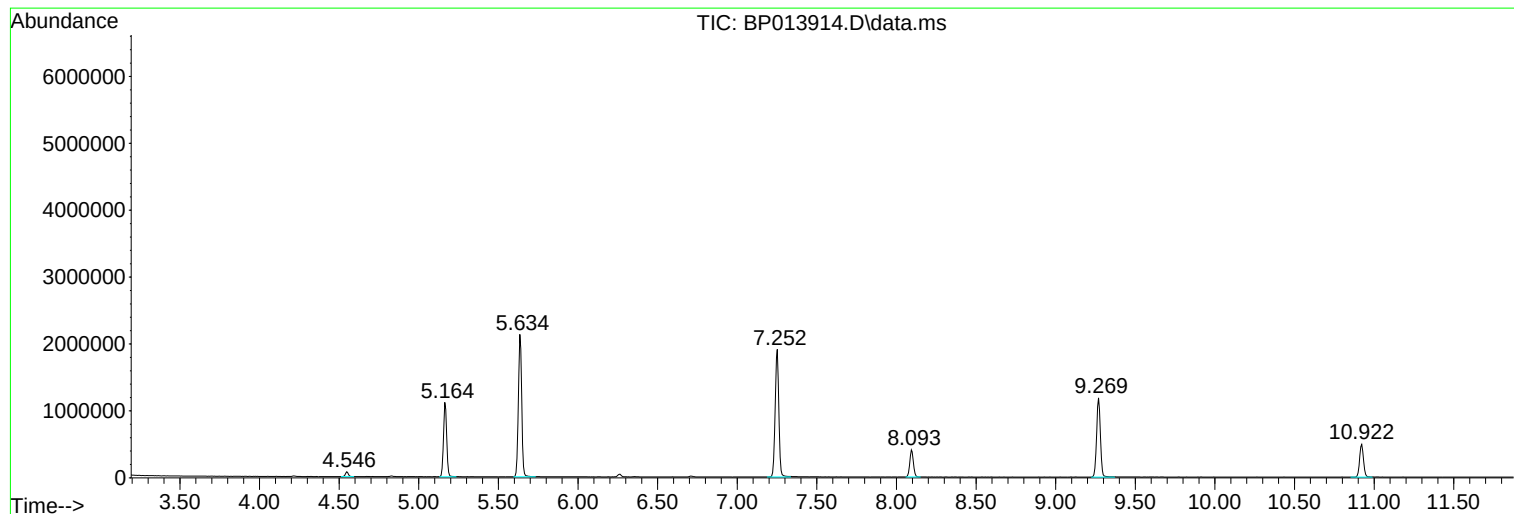
Sum of corrected areas: 33749625

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP021623.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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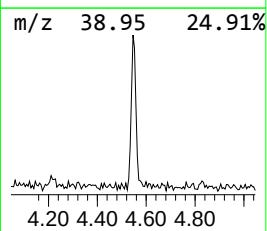
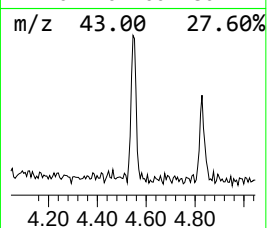
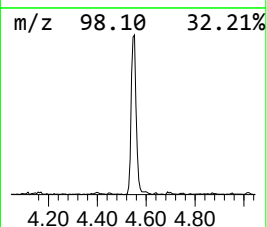
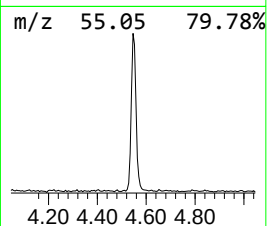
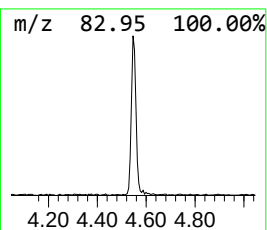
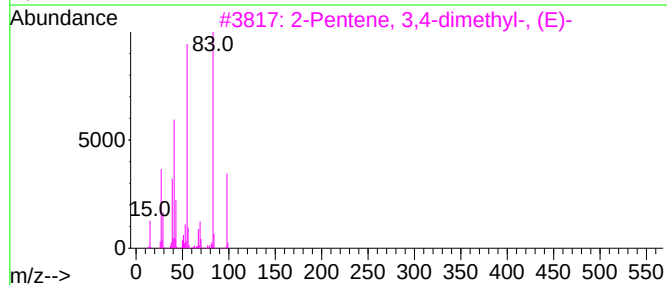
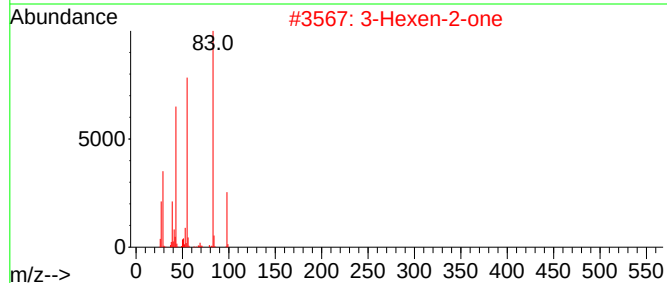
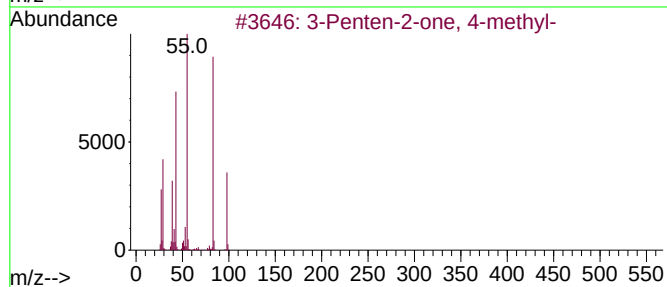
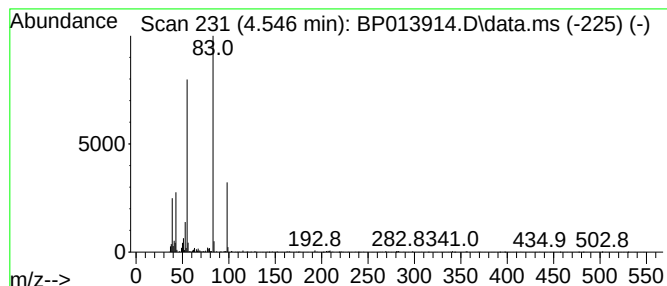
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TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.546	3.03 ng	96900	1,4-Dichlorobenzene-d4	8.093

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	87
3			2-Pentene, 3,4-dimethyl-, (E)-	98	C7H14	004914-92-5	86
4			2-Pentene, 3,4-dimethyl-, (Z)-	98	C7H14	004914-91-4	86
5			Cyclohexane, methyl-	98	C7H14	000108-87-2	80



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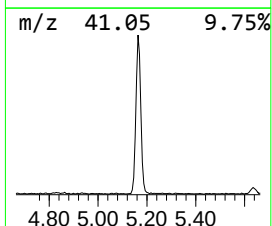
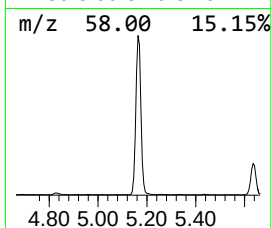
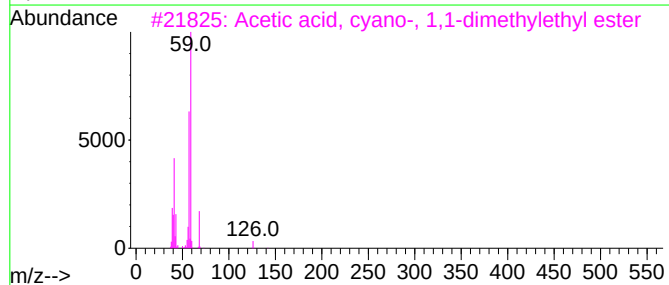
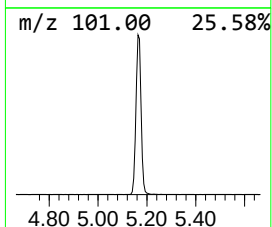
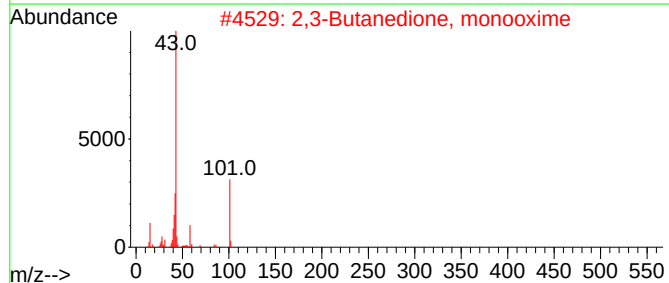
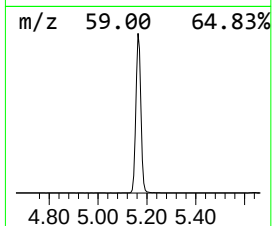
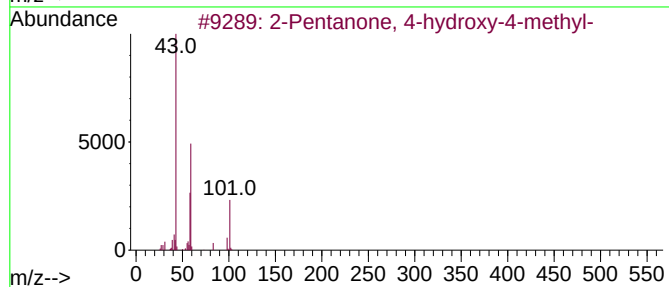
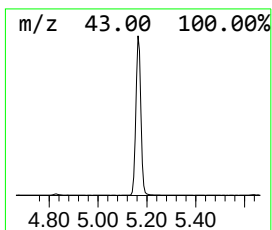
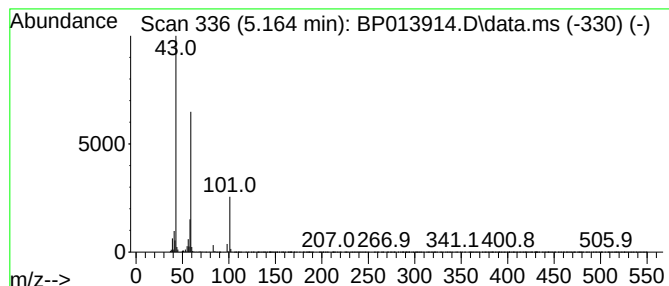
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Peak Number 2 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.164	48.88 ng	1563530	1,4-Dichlorobenzene-d4	8.093

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56
2			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	25
3			Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	17
4			1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	10
5			5-Hexen-2-one	98	C6H10O	000109-49-9	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
3-Penten-2-one,...	4.546	3.0	ng	96900	1	8.093	639713	20.0
2-Pentanone, 4-...	5.164	48.9	ng	1563530	1	8.093	639713	20.0