

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG111722\
 Data File : BG055659.D
 Acq On : 17 Nov 2022 10:07
 Operator : CG/JU
 Sample : PB149062BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB149062BL

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_G\Methods\8270-BG111622.M
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BG055659.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	4.685	232	238	246	rVB	30238	46778	1.63%	0.420%
2	5.302	337	343	353	rVB	186081	286822	9.98%	2.573%
3	5.778	417	424	437	rBV	569836	915504	31.87%	8.213%
4	7.394	690	699	707	rBV	569461	969581	33.75%	8.698%
5	8.252	837	845	853	rBV	97750	164685	5.73%	1.477%
6	9.421	1036	1044	1054	rVB	330325	588587	20.49%	5.280%
7	11.084	1319	1327	1336	rBV	138902	248039	8.63%	2.225%
8	13.499	1730	1738	1746	rBV	936045	1469024	51.14%	13.179%
9	14.873	1966	1972	1982	rVB	239188	375808	13.08%	3.371%
10	16.354	2217	2224	2243	rBV	857522	1334003	46.44%	11.968%
11	17.617	2432	2439	2446	rBV	351325	537367	18.71%	4.821%
12	20.202	2873	2879	2885	rBV	2185446	2872742	100.00%	25.772%
13	21.912	3164	3170	3180	rVB2	384761	646820	22.52%	5.803%
14	25.332	3741	3752	3763	rVB2	234174	690959	24.05%	6.199%

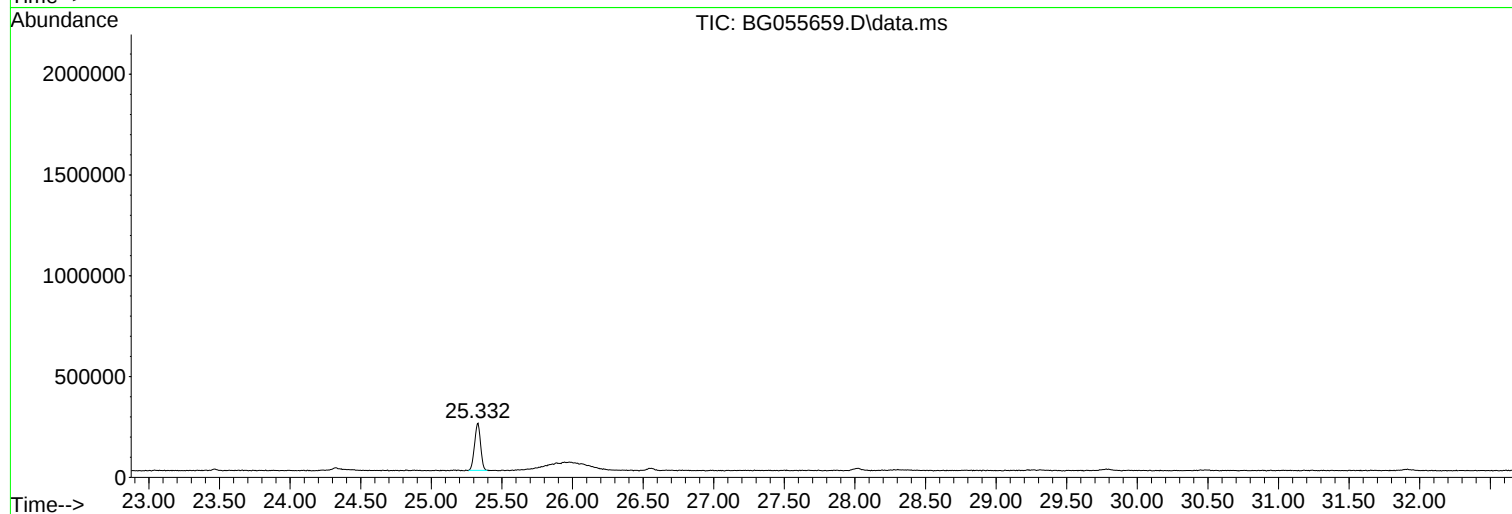
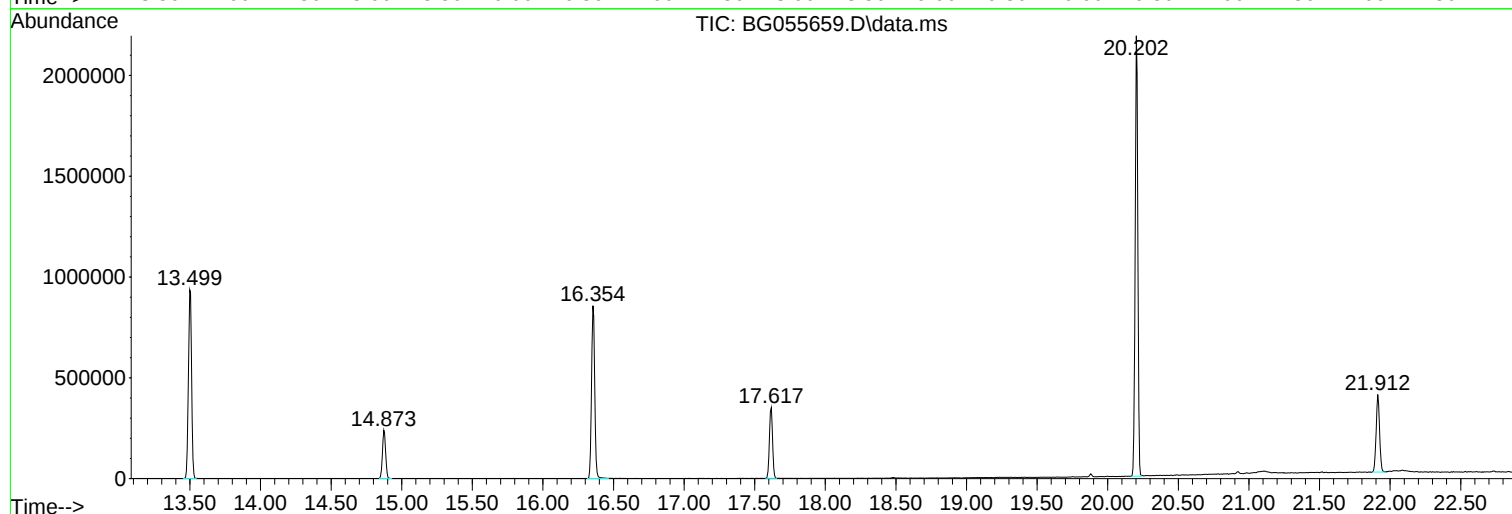
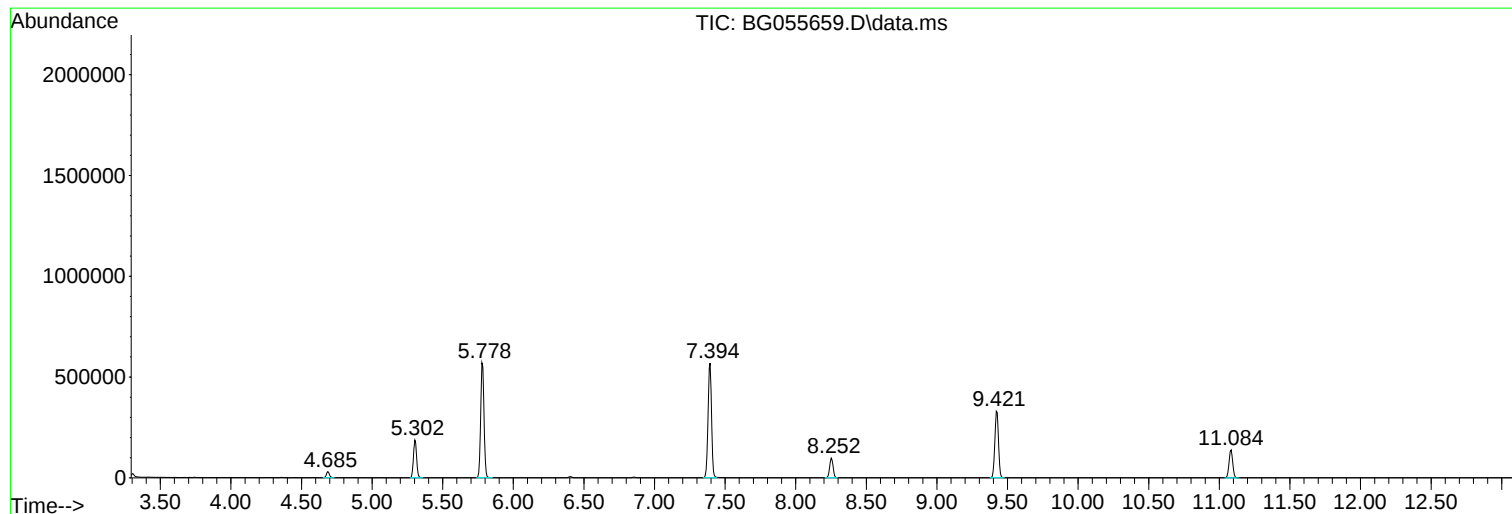
Sum of corrected areas: 11146719

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG111622.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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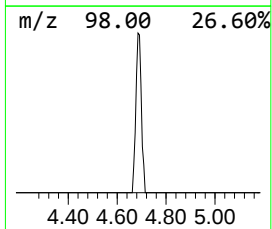
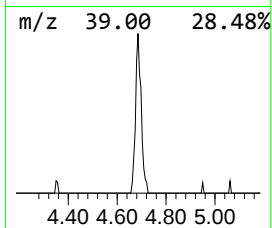
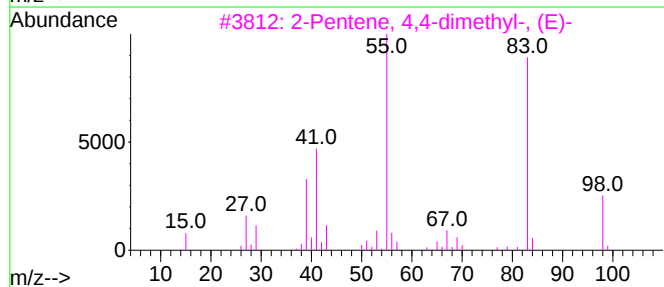
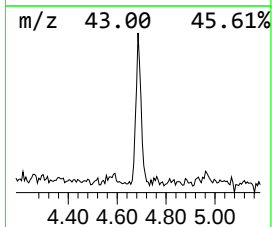
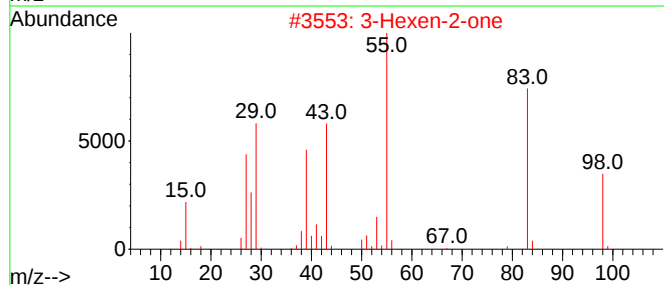
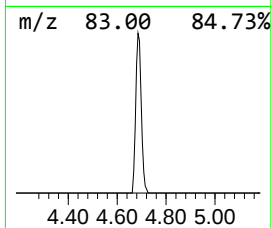
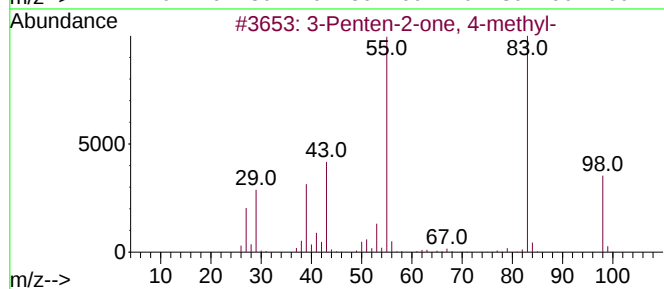
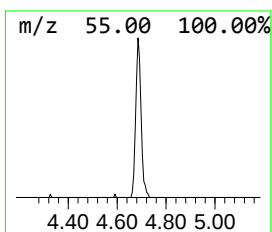
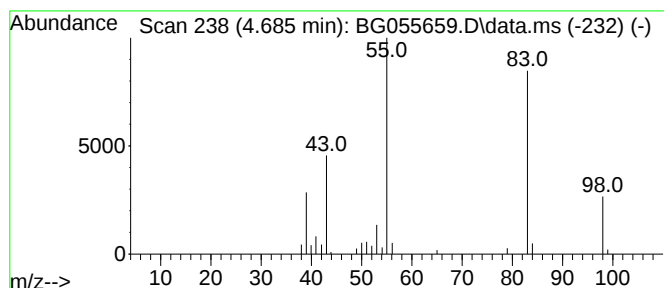
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG111622.M
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.685	5.68 ng	46778	1,4-Dichlorobenzene-d4	8.252

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, 4,4-dimethyl-, (E)-	98	C7H14	000690-08-4	86
4			2-Pentene, 2,4-dimethyl-	98	C7H14	000625-65-0	86
5			2-Pentene, 3,4-dimethyl-	98	C7H14	024910-63-2	86



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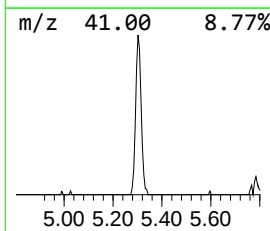
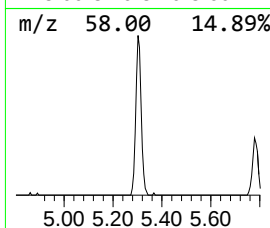
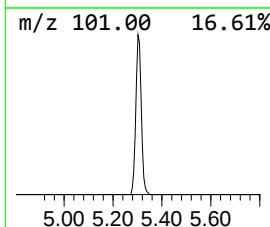
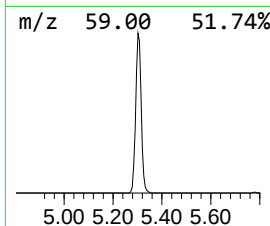
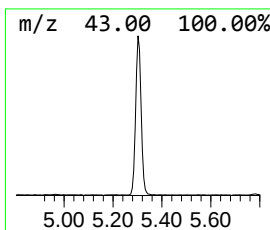
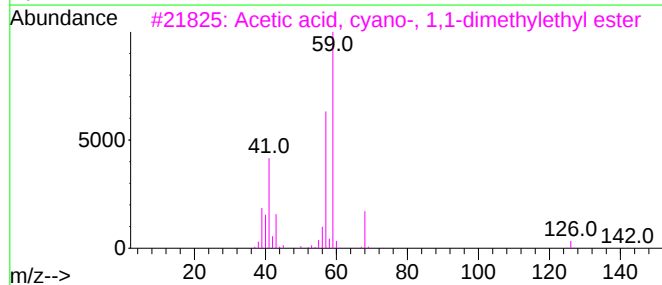
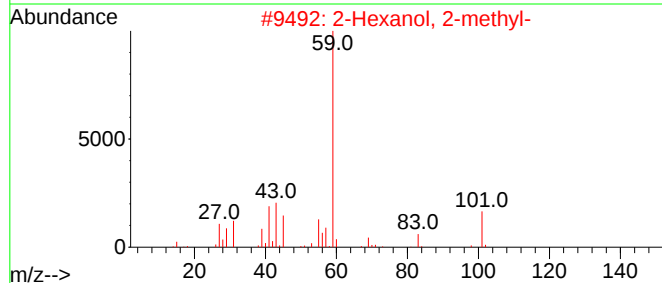
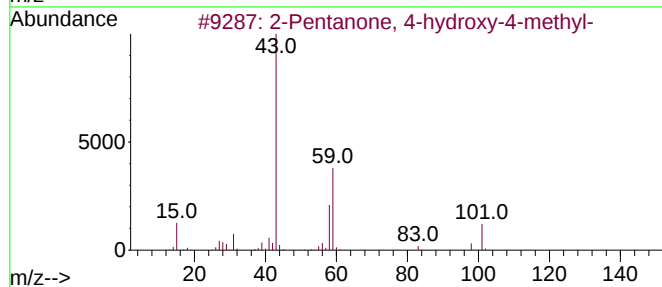
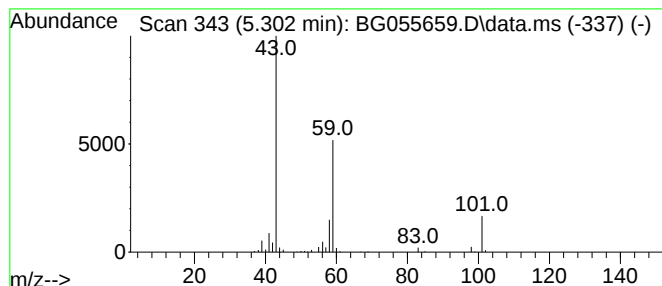
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Peak Number 2 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.302	34.83 ng	286822	1,4-Dichlorobenzene-d4	8.252

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	59
2		2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	28
3		Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	25
4		2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	16
5		1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	12



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
3-Penten-2-one,...	4.685	5.7	ng	46778	1	8.252	164685	20.0
2-Pentanone, 4-...	5.302	34.8	ng	286822	1	8.252	164685	20.0