

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG122122\
 Data File : BG056043.D
 Acq On : 21 Dec 2022 10:56
 Operator : CG/JU
 Sample : PB149688BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB149688BL

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-BG121322.M
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BG056043.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.400	354	359	365	rVB	14842	21336	1.80%	0.578%
2	5.881	433	441	448	rBB	129610	207063	17.43%	5.610%
3	7.491	707	715	722	rBB	134282	226188	19.04%	6.128%
4	8.361	857	863	869	rBB	25913	42858	3.61%	1.161%
5	9.536	1055	1063	1069	rBB	78345	140321	11.81%	3.801%
6	11.193	1339	1345	1352	rBB	34950	59367	5.00%	1.608%
7	13.596	1748	1754	1762	rBB	306314	463277	39.00%	12.551%
8	14.971	1981	1988	1994	rBV	73747	110109	9.27%	2.983%
9	16.446	2233	2239	2246	rBV	378595	561674	47.29%	15.216%
10	17.709	2447	2454	2462	rBB	117666	177752	14.96%	4.816%
11	20.282	2886	2892	2897	rBV	906384	1187809	100.00%	32.179%
12	22.010	3179	3186	3195	rVB	141627	238827	20.11%	6.470%
13	25.494	3769	3779	3794	rVB2	81786	254640	21.44%	6.899%

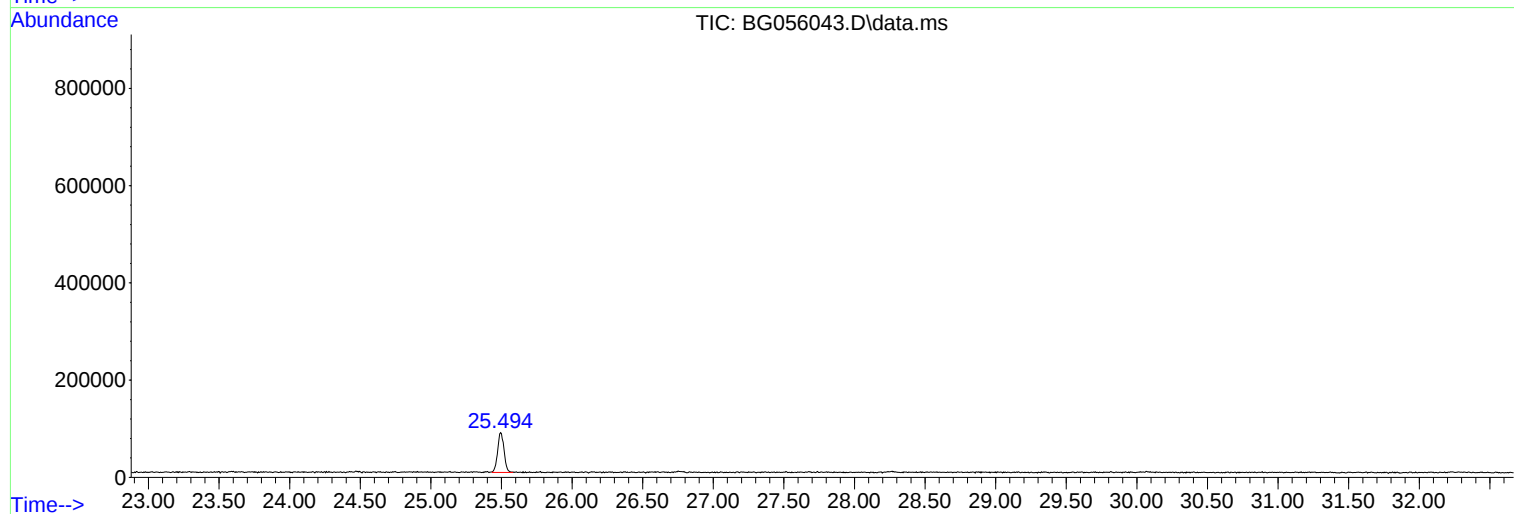
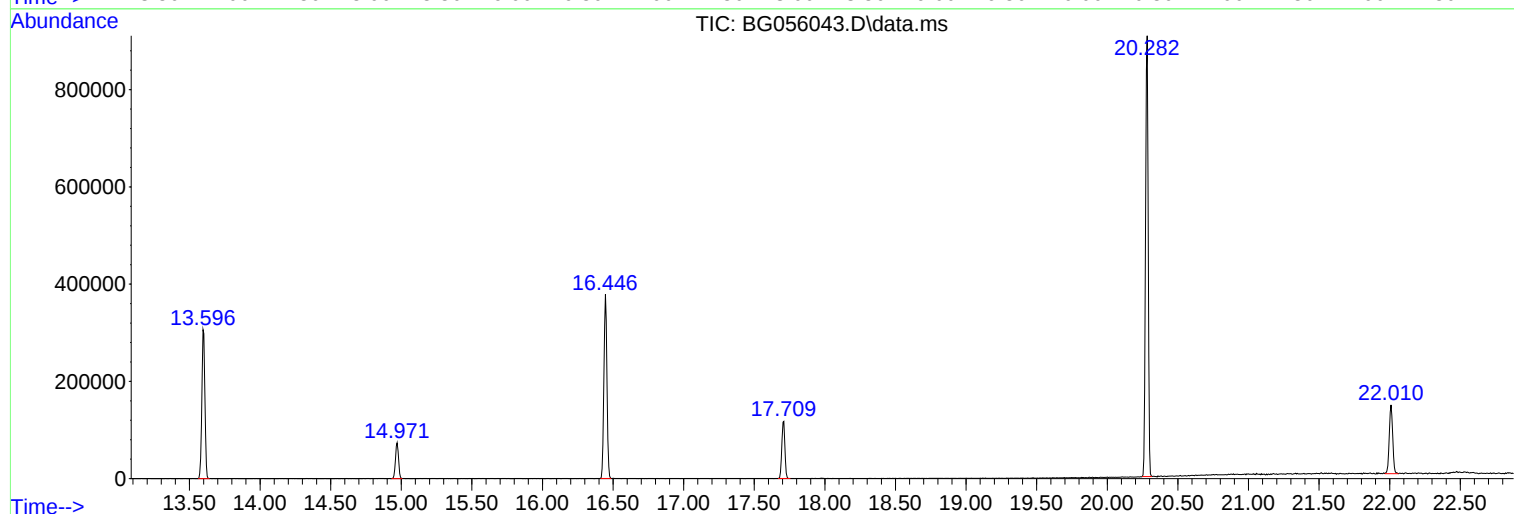
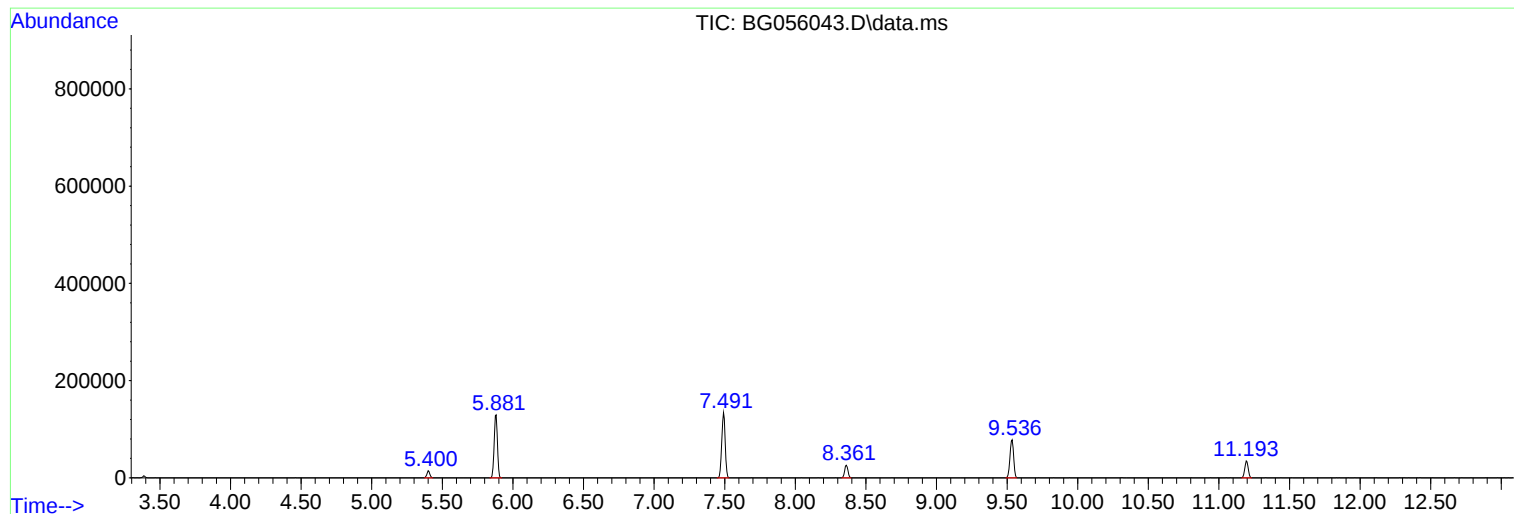
Sum of corrected areas: 3691221

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG121322.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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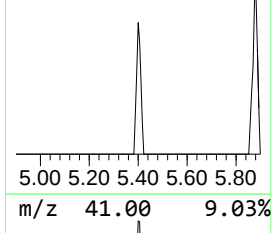
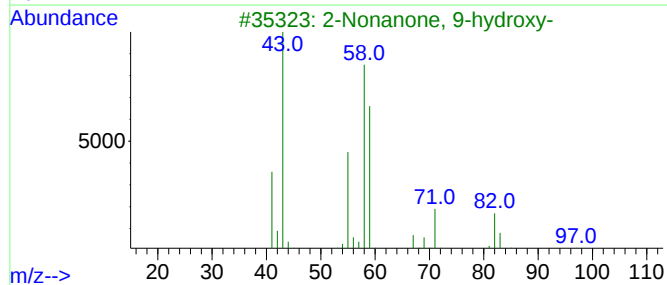
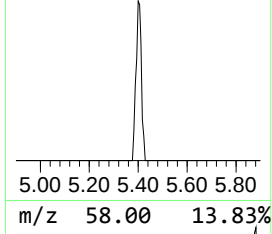
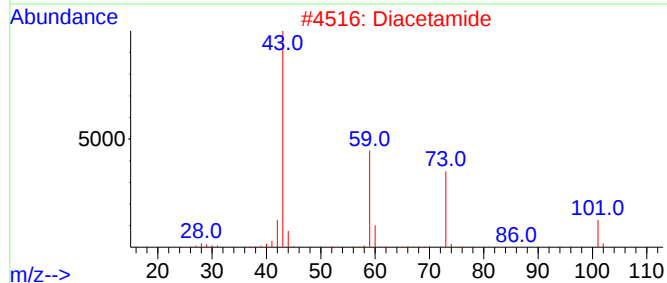
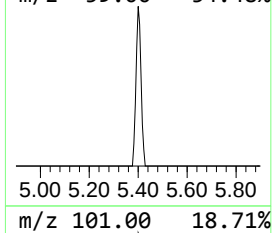
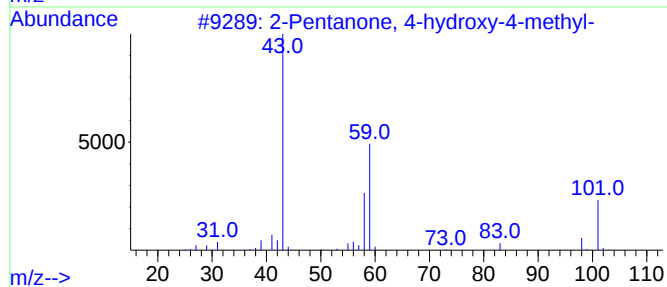
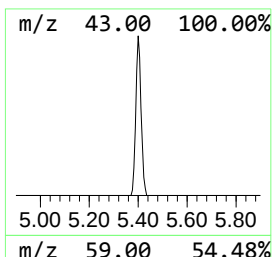
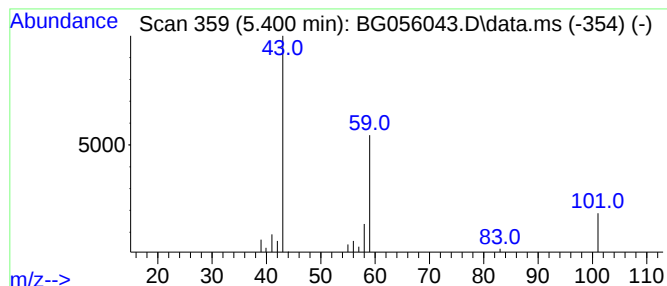
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.400	9.96 ng	21336	1,4-Dichlorobenzene-d4	8.361

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	72		
2	Diacetamide	101	C4H7NO2	000625-77-4	9		
3	2-Nonanone, 9-hydroxy-	158	C9H18O2	025368-56-3	9		
4	1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	9		
5	2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	9		



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-...	5.400	10.0	ng	21336	1	8.361	42858	20.0