

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG102022\
 Data File : BG055218.D
 Acq On : 20 Oct 2022 11:58
 Operator : CG/JU
 Sample : PB148413BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB148413BL

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-BG101922.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BG055218.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.321	340	346	358	rVB	153387	230748	6.64%	1.643%
2	5.803	421	428	441	rBV	853481	1317296	37.92%	9.379%
3	7.419	695	703	717	rBV	813652	1428886	41.13%	10.174%
4	8.283	843	850	857	rBV	128888	217419	6.26%	1.548%
5	9.458	1041	1050	1063	rBV	476470	847578	24.40%	6.035%
6	11.115	1324	1332	1346	rBV	184827	345474	9.94%	2.460%
7	13.529	1736	1743	1753	rBV	1319547	1994374	57.41%	14.200%
8	14.898	1966	1976	1987	rBV	336633	510979	14.71%	3.638%
9	16.373	2220	2227	2245	rBV	1029097	1582599	45.56%	11.269%
10	17.631	2434	2441	2450	rBV	454436	675461	19.44%	4.809%
11	20.204	2873	2879	2890	rBV	2543316	3473946	100.00%	24.735%
12	21.908	3163	3169	3178	rBV2	419719	706606	20.34%	5.031%
13	25.322	3739	3750	3763	rBV	234069	713054	20.53%	5.077%

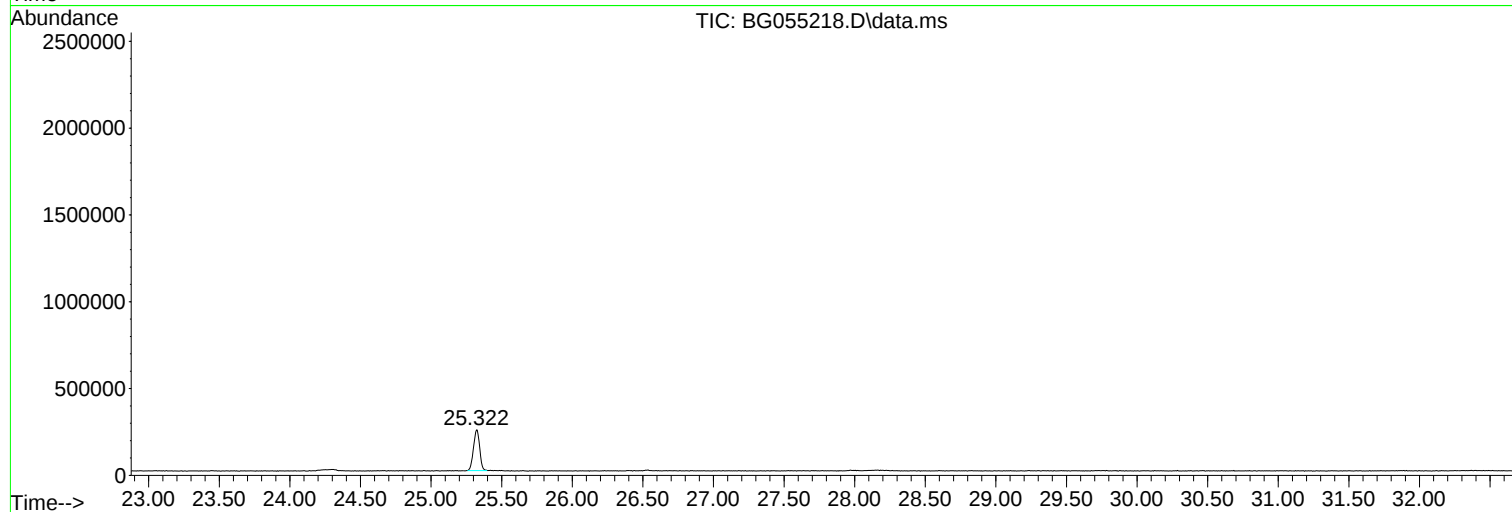
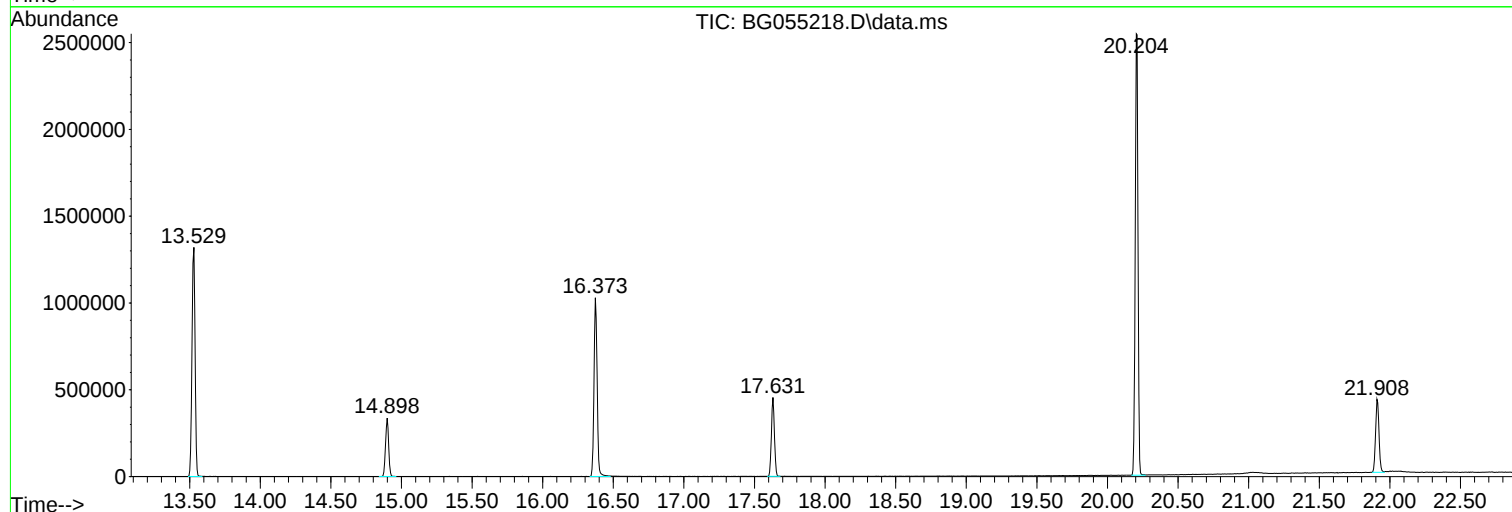
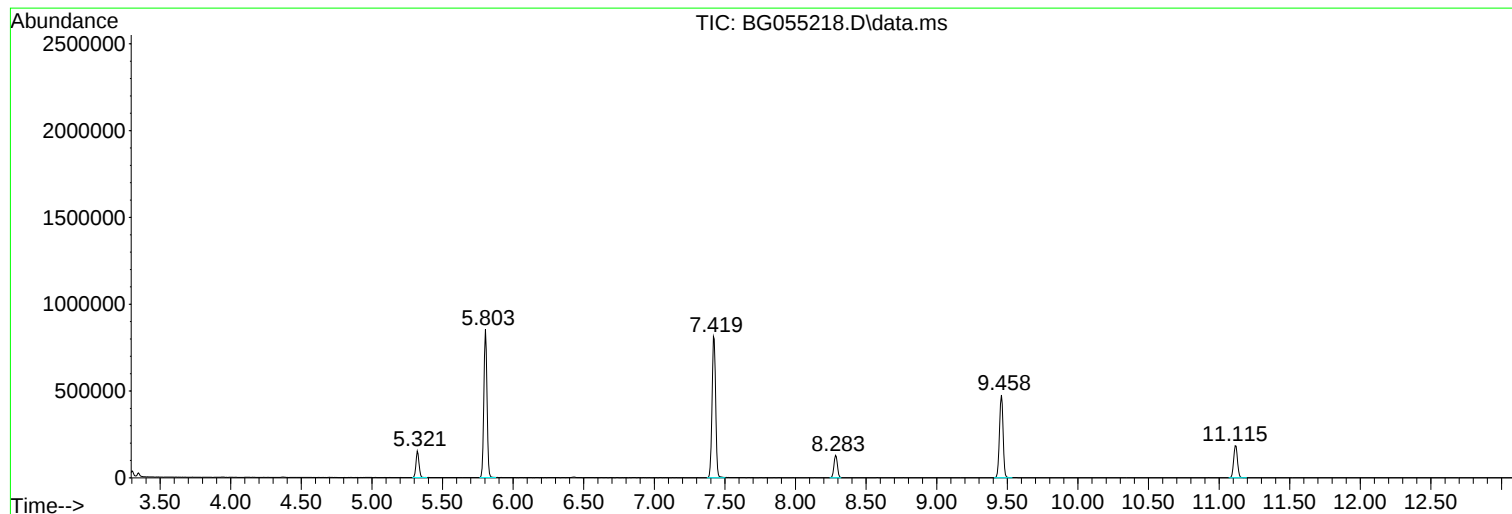
Sum of corrected areas: 14044420

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG101922.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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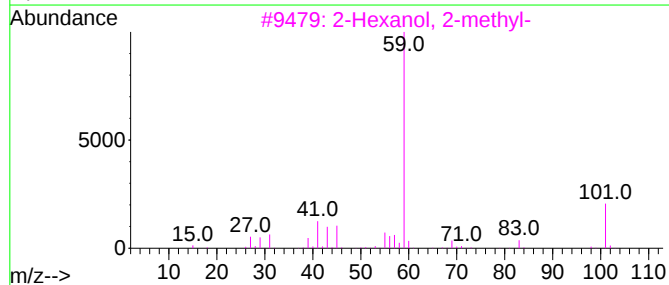
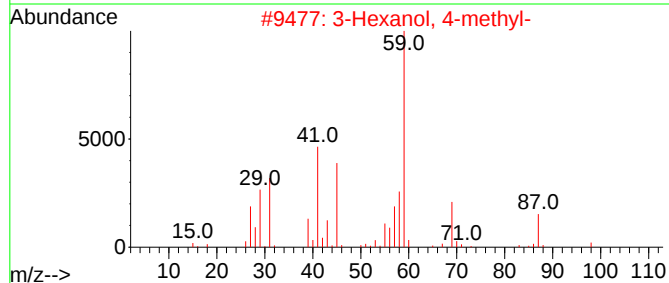
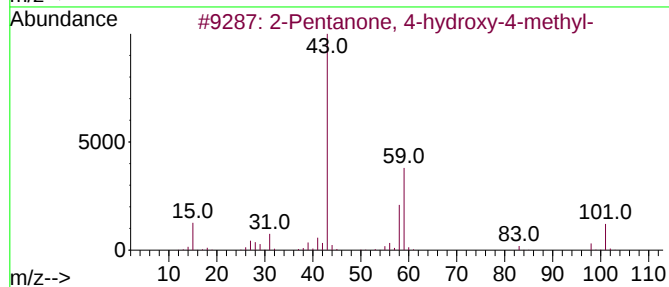
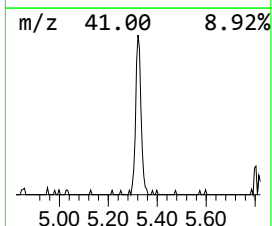
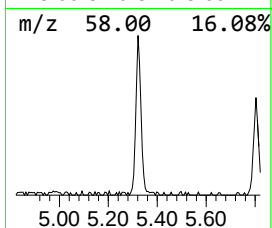
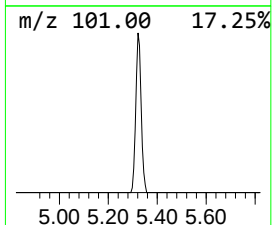
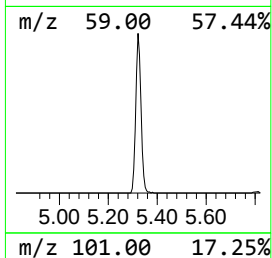
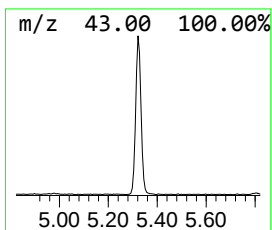
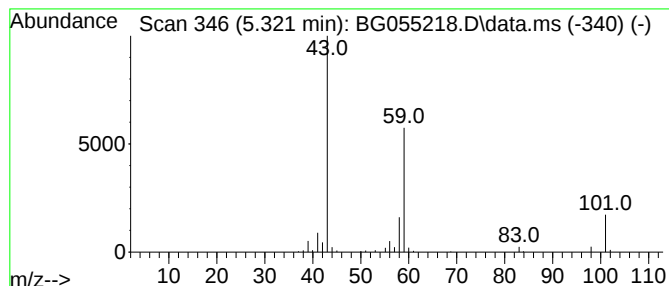
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.321	21.23 ng	230748	1,4-Dichlorobenzene-d4	8.283

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	59		
2	3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	33		
3	2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	28		
4	Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	25		
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
2-Pentanone, 4-...	5.321	21.2	ng	230748	1	8.283	217419	20.0