

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM081224\
Data File : BM047153.D
Acq On : 12 Aug 2024 12:02
Operator : RC/JU
Sample : PB162517BL
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
PB162517BL

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_M\Methods\8270-BM081024.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BM047153.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	4.581	283	288	295	rBV	292690	398449	3.89%	0.809%
2	5.028	357	364	378	rBV	3578511	5143123	50.26%	10.440%
3	6.587	621	629	642	rBV	3447097	5424086	53.00%	11.010%
4	7.375	755	763	769	rBV	703360	1088220	10.63%	2.209%
5	8.522	949	958	969	rBV	2107328	3506421	34.26%	7.117%
6	10.128	1224	1231	1248	rVB	914929	1472716	14.39%	2.989%
7	12.639	1651	1658	1674	rBV	4896578	7185327	70.21%	14.585%
8	14.027	1888	1894	1909	rBV	1335670	1925970	18.82%	3.909%
9	15.533	2144	2150	2165	rBV	4281749	5963250	58.27%	12.104%
10	16.792	2357	2364	2376	rBV	1646367	2289457	22.37%	4.647%
11	19.456	2811	2817	2825	rBV	8365694	10233791	100.00%	20.773%
12	21.027	3079	3084	3097	rBV	1628112	2286279	22.34%	4.641%
13	23.721	3535	3542	3553	rVB	1053895	2348685	22.95%	4.767%

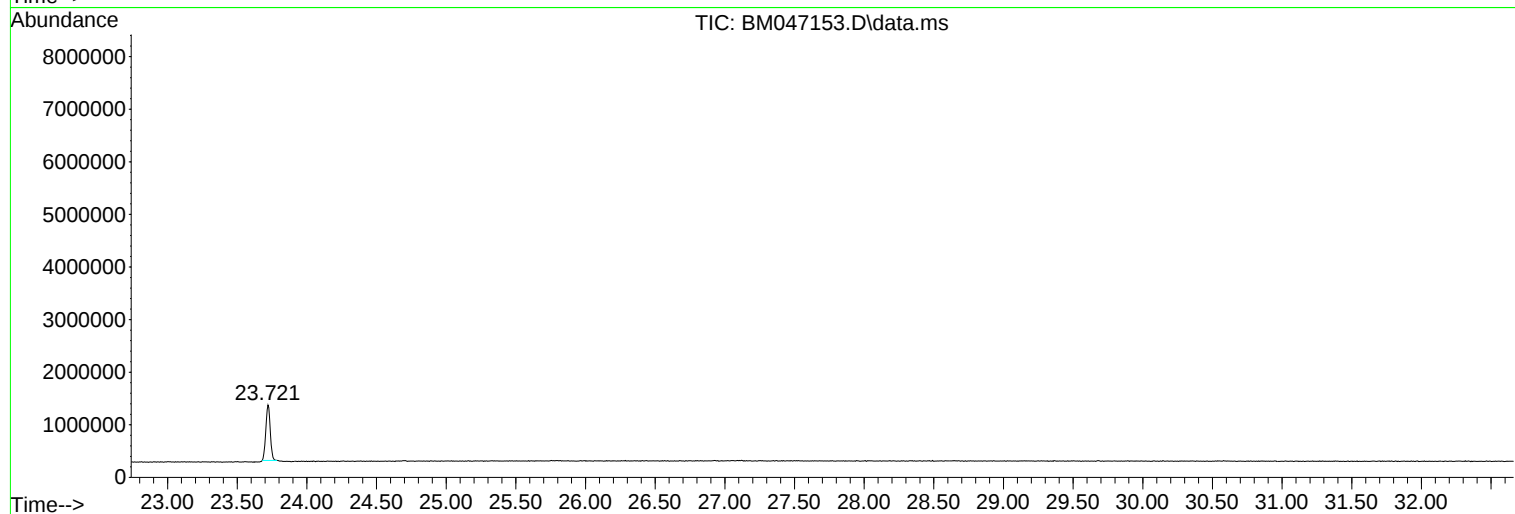
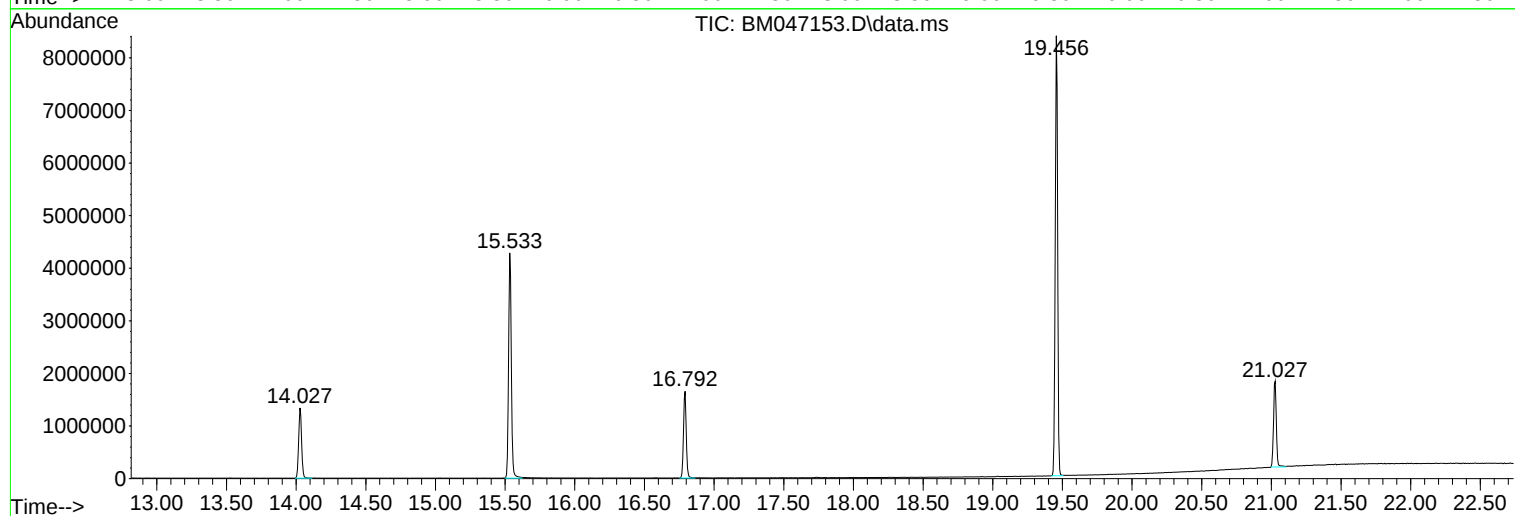
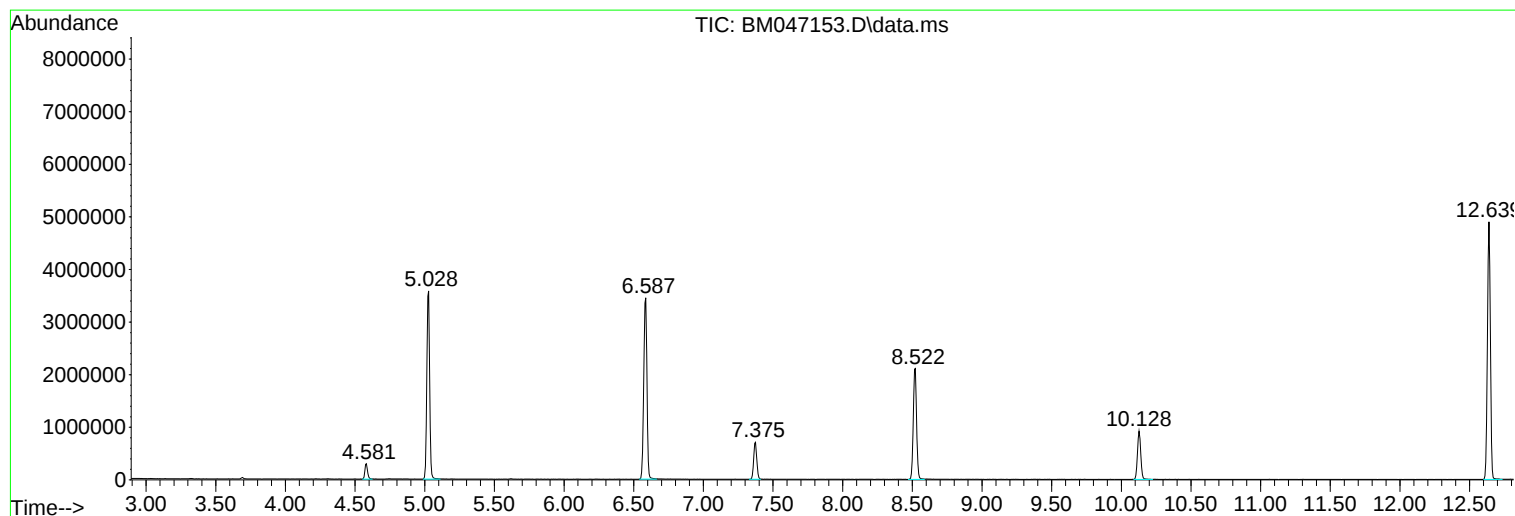
Sum of corrected areas: 49265774

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM081024.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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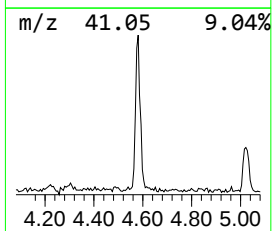
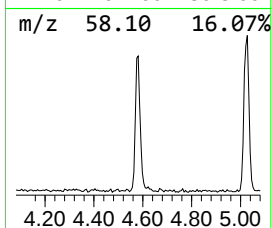
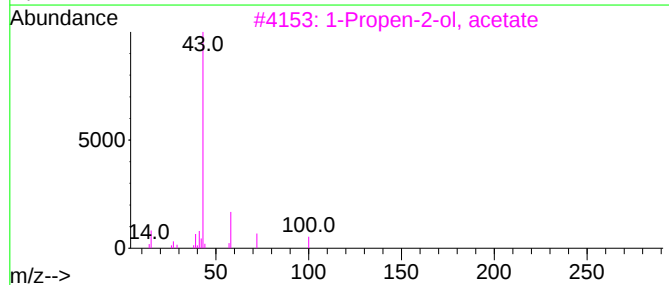
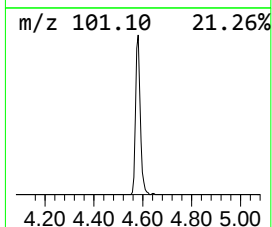
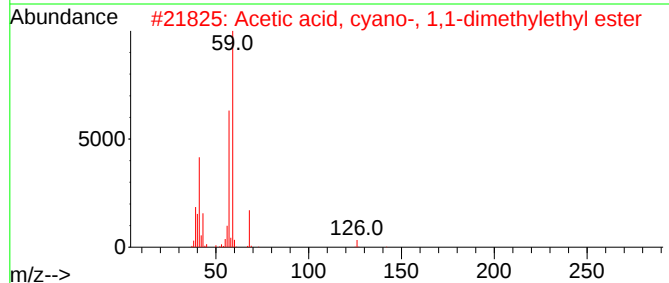
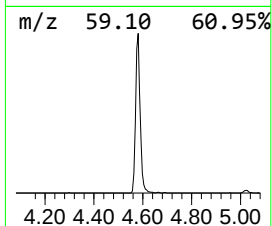
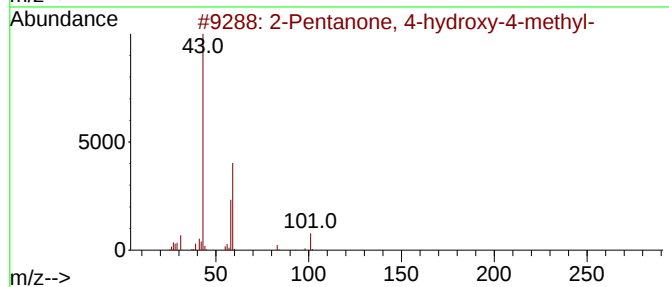
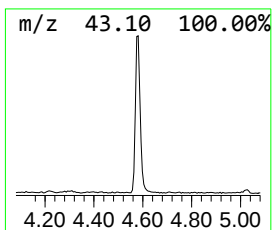
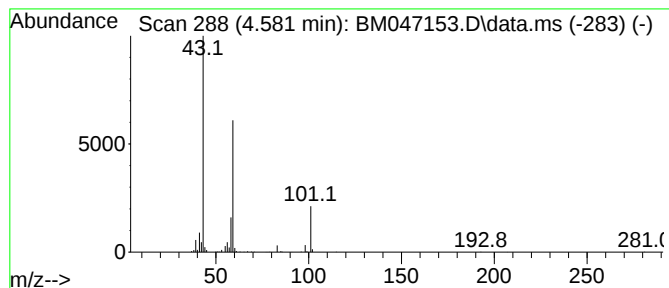
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.581	7.32 ng	398449	1,4-Dichlorobenzene-d4	7.375

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56		
2	Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	23		
3	1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	10		
4	5-Hexen-2-one	98	C6H10O	000109-49-9	9		
5	Butane, 1-ethoxy-	102	C6H14O	000628-81-9	9		



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-...	4.581	7.3	ng	398449	1	7.375	1088220	20.0