

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM103020\
 Data File : BM027666.D
 Acq On : 30 Oct 2020 19:02
 Operator : JU/CG
 Sample : PB132643BL
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 PB132643BL

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-BM103020.M
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BM027666.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.381	401	407	414	rBV	79016	114989	5.88%	1.184%
2	5.875	484	491	498	rBV	764663	1151587	58.86%	11.861%
3	7.522	763	771	778	rBV	764144	1210202	61.86%	12.465%
4	8.404	914	921	928	rBV	135988	214131	10.94%	2.205%
5	9.581	1111	1121	1128	rBV	492986	822262	42.03%	8.469%
6	11.245	1397	1404	1412	rBV	168620	275362	14.07%	2.836%
7	13.651	1806	1813	1819	rBV	1072503	1513236	77.35%	15.586%
8	15.016	2038	2045	2051	rVB	226985	332392	16.99%	3.424%
9	16.474	2287	2293	2302	rBV	714978	1007757	51.51%	10.380%
10	17.727	2500	2506	2513	rVB	264686	371933	19.01%	3.831%
11	20.286	2935	2941	2946	rVB	1592766	1956470	100.00%	20.151%
12	21.821	3197	3202	3209	rVB	286268	375871	19.21%	3.871%
13	24.280	3612	3620	3628	rVB2	178584	362913	18.55%	3.738%

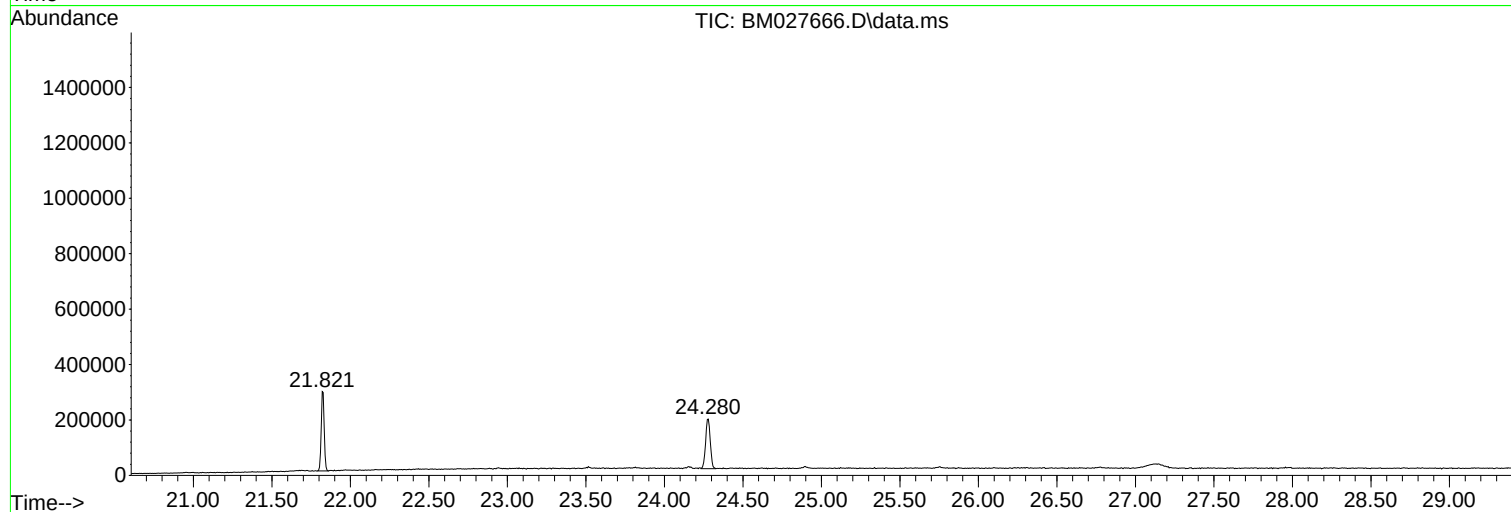
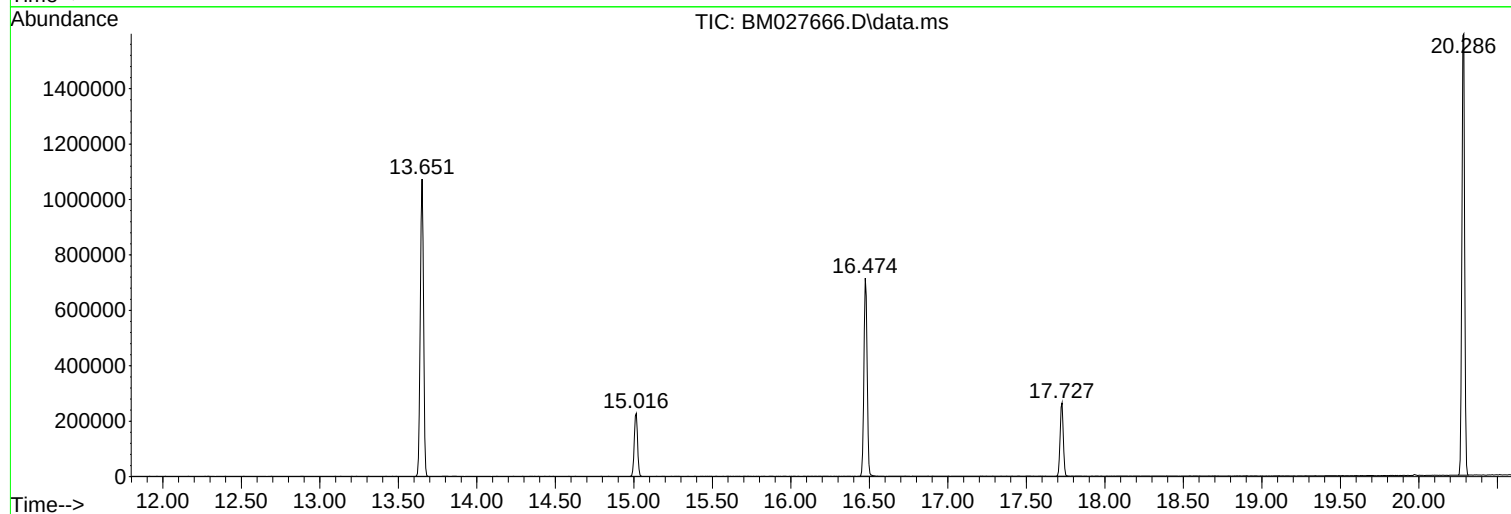
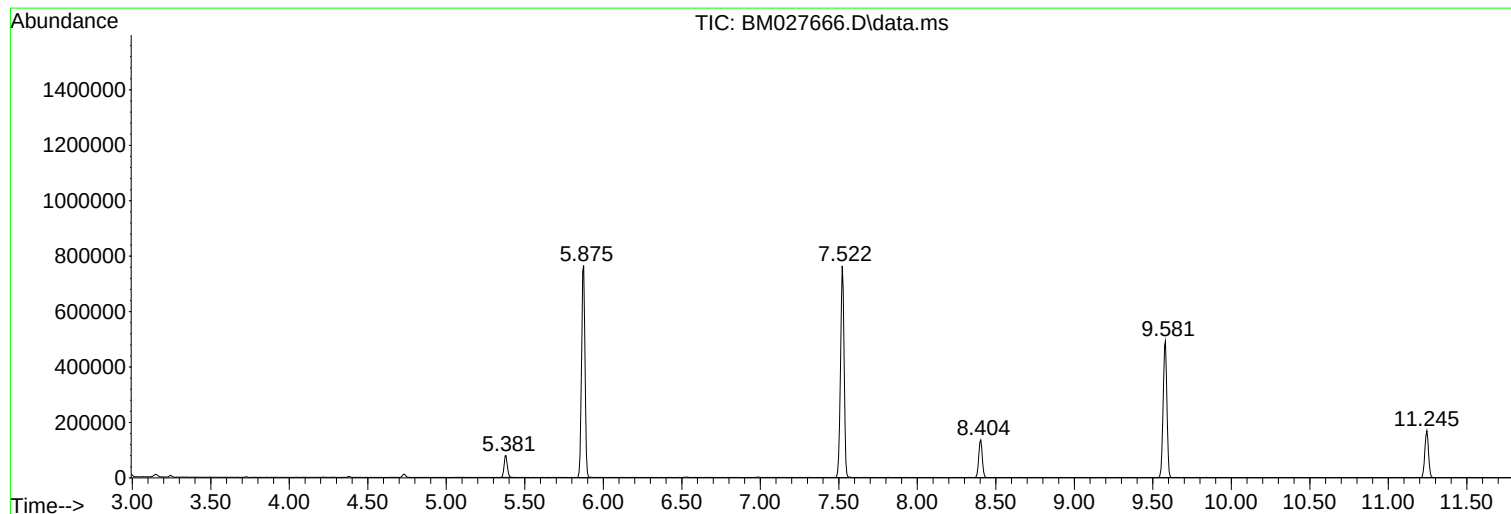
Sum of corrected areas: 9709105

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM103020.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P



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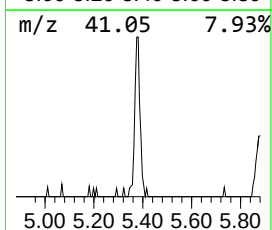
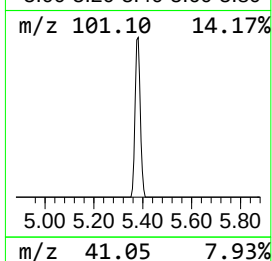
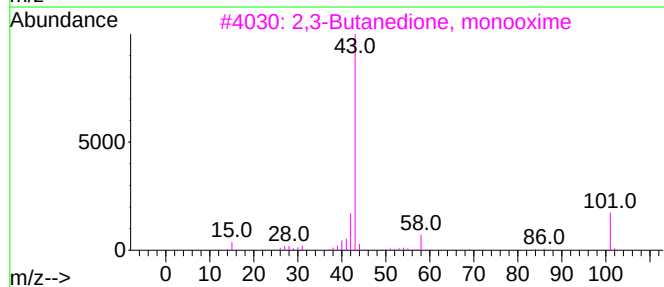
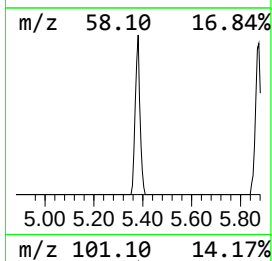
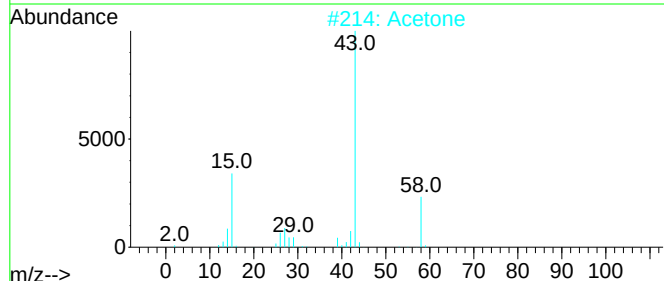
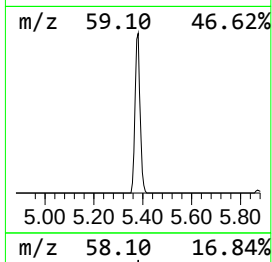
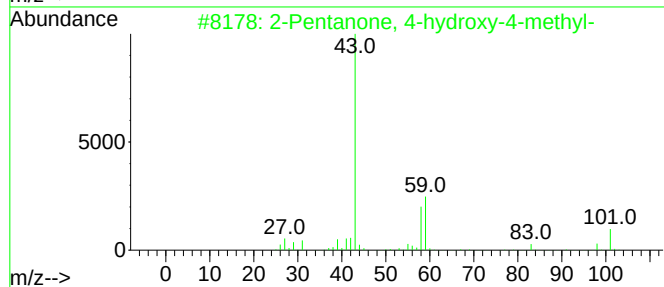
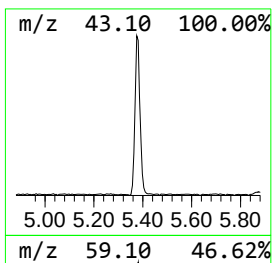
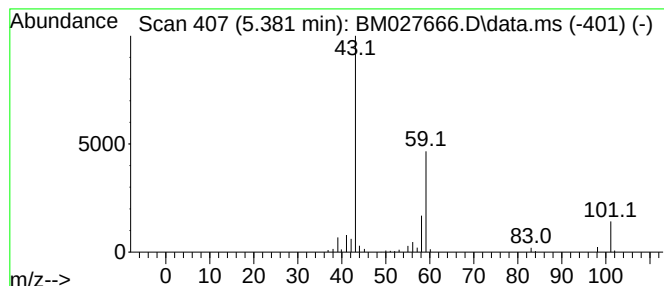
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.381	10.74 ng	114989	1,4-Dichlorobenzene-d4	8.404

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	64		
2	Acetone	58	C3H6O	000067-64-1	12		
3	2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	10		
4	Diacetamide	101	C4H7NO2	000625-77-4	9		
5	5-Oxohexanenitrile	111	C6H9NO	010412-98-3	9		



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-...	5.381	10.7	ng	114989	1	8.404	214131	20.0