

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM061221\
 Data File : BM030413.D
 Acq On : 12 Jun 2021 13:59
 Operator : CG/JU
 Sample : PB137009BL
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 PB137009BL

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-BM060921.M
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BM030413.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	4.963	314	319	331	rVB	163214	258618	1.91%	0.497%
2	5.416	389	396	411	rBV	3301114	4915171	36.38%	9.444%
3	6.998	655	665	679	rBV	3307492	5182479	38.36%	9.957%
4	7.834	799	807	814	rBV	544131	860619	6.37%	1.654%
5	8.986	995	1003	1021	rBV	1729338	2952940	21.86%	5.674%
6	10.622	1273	1281	1290	rBV	710205	1208476	8.94%	2.322%
7	13.080	1691	1699	1718	rBV	5218971	7555542	55.92%	14.517%
8	14.457	1926	1933	1946	rBV	1145982	1692335	12.53%	3.252%
9	15.945	2179	2186	2207	rBV	3947111	5945212	44.00%	11.423%
10	17.192	2392	2398	2412	rBV	1473479	2138663	15.83%	4.109%
11	19.809	2837	2843	2848	rBV	11018899	13510622	100.00%	25.959%
12	21.356	3101	3106	3115	rBV	2190901	2782546	20.60%	5.346%
13	23.633	3487	3493	3505	rVB2	1592177	3042953	22.52%	5.847%

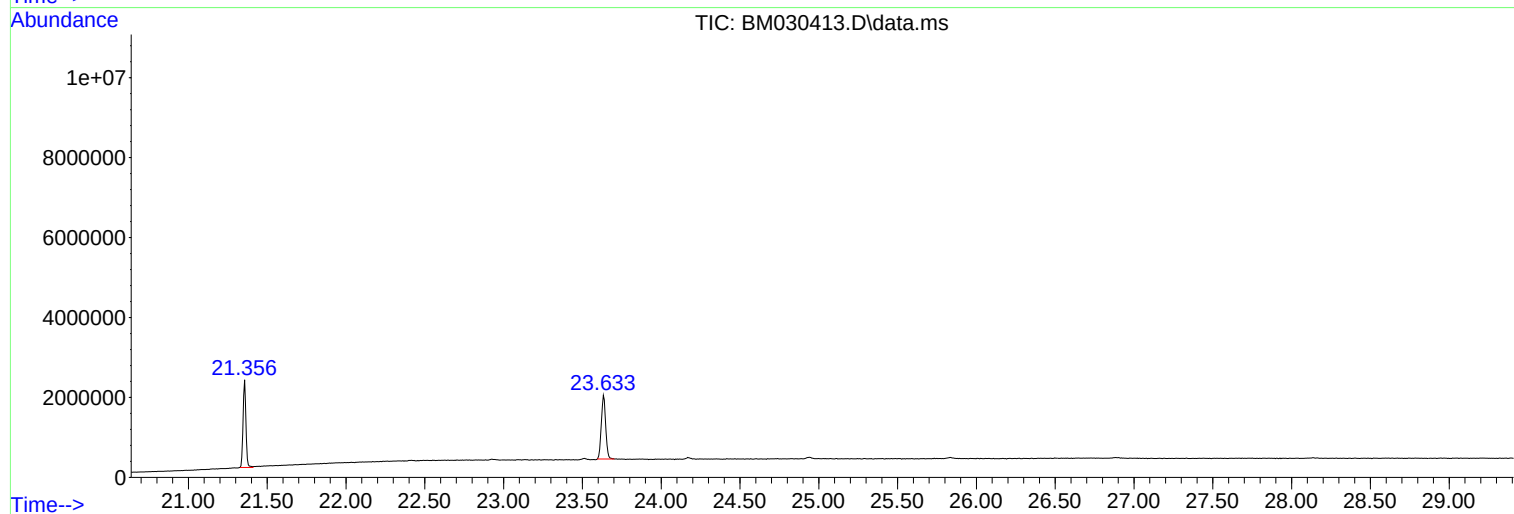
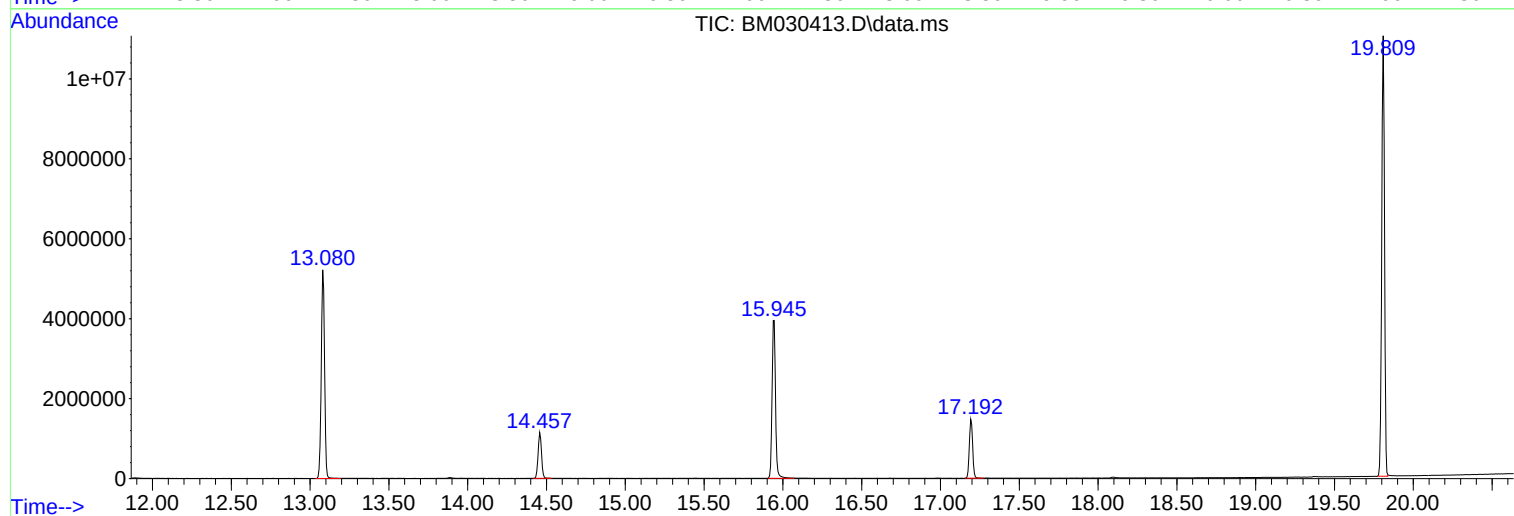
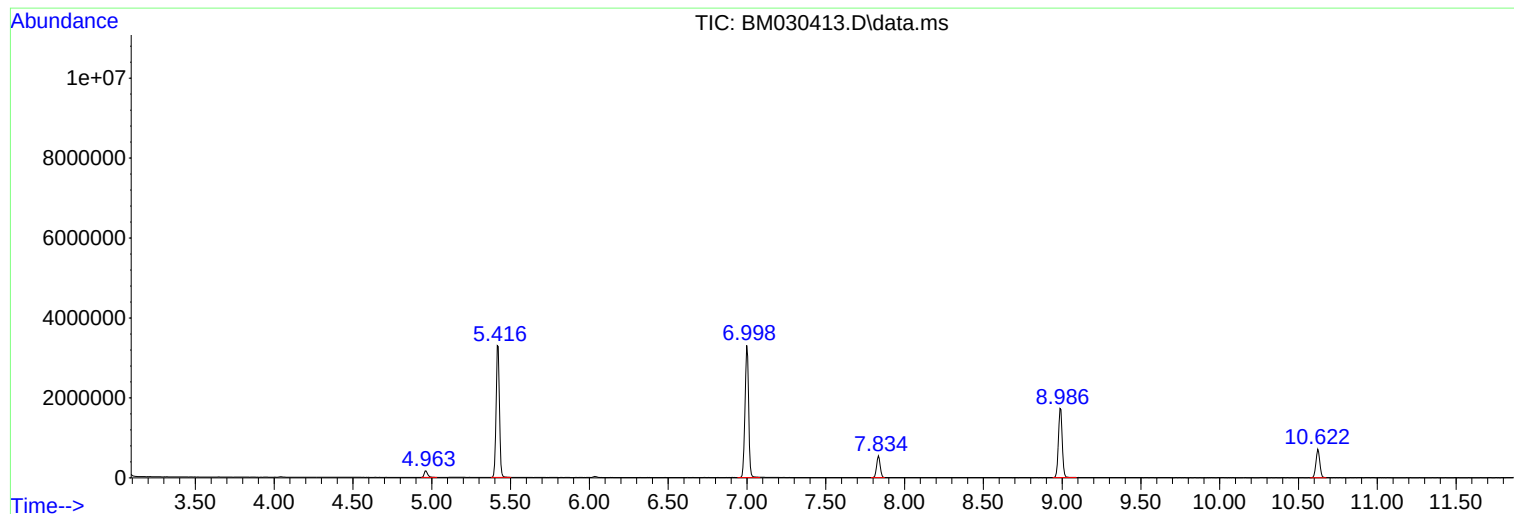
Sum of corrected areas: 52046176

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM060921.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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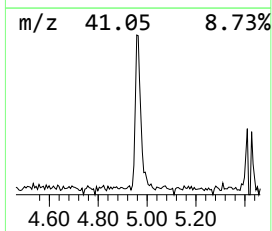
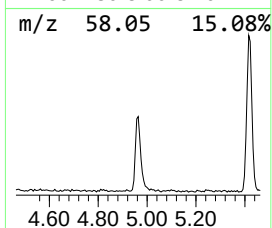
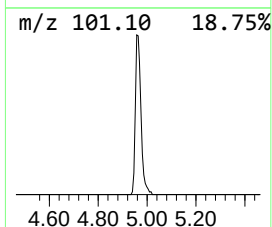
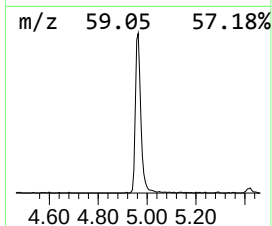
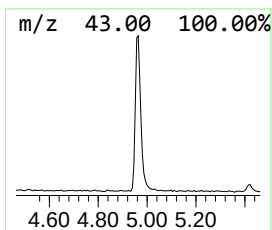
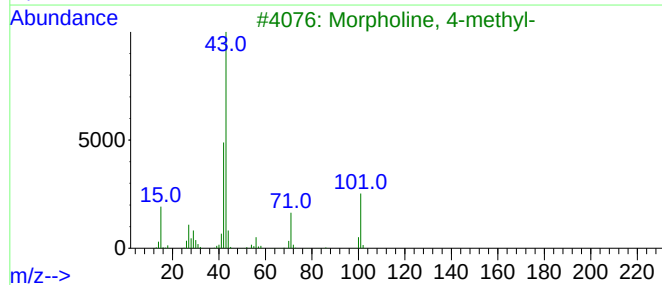
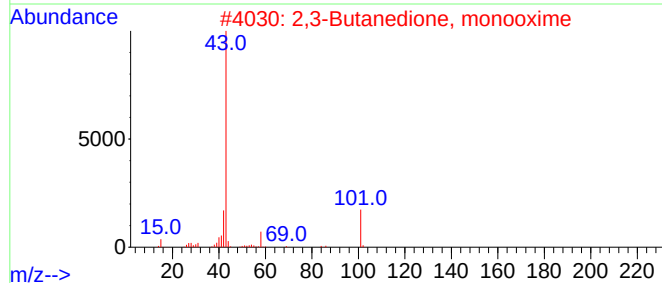
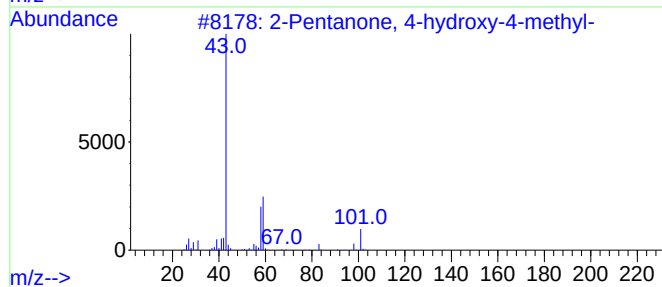
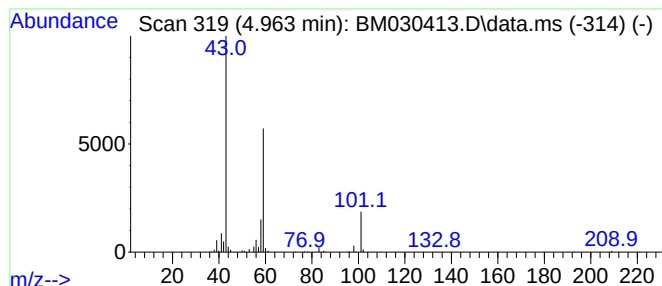
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.963	6.01 ng	258618	1,4-Dichlorobenzene-d4	7.834

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50		
2	2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	16		
3	Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9		
4	Butane, 1-ethoxy-	102	C6H14O	000628-81-9	9		
5	(+)-4-Amino-4,5-dihydro-2(3H)-f...	101	C4H7NO2	016504-58-8	9		



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-...	4.963	6.0	ng	258618	1	7.834	860619	20.0