

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM061221\
 Data File : BM030417.D
 Acq On : 12 Jun 2021 16:32
 Operator : CG/JU
 Sample : M2673-02
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 17-B6(92-96)

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: BM030417.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	4.969	315	320	333	rVB	68487	109472	1.46%	0.310%
2	5.416	389	396	409	rBV	2088934	3004607	40.08%	8.515%
3	6.998	657	665	682	rBV	1952869	3096693	41.31%	8.776%
4	7.834	800	807	815	rBV	594689	937067	12.50%	2.656%
5	8.986	994	1003	1014	rBV	1100582	1832958	24.45%	5.194%
6	10.422	1239	1247	1262	rBV	31202	93586	1.25%	0.265%
7	10.622	1273	1281	1291	rVB	737257	1242148	16.57%	3.520%
8	13.080	1692	1699	1713	rBV	2992173	4379880	58.43%	12.412%
9	14.457	1926	1933	1947	rVB	1101363	1639111	21.87%	4.645%
10	15.939	2175	2185	2204	rBV	2306815	3307561	44.13%	9.373%
11	17.192	2392	2398	2410	rBV	1368536	1993689	26.60%	5.650%
12	18.080	2545	2549	2558	rBV	105754	188474	2.51%	0.534%
13	19.356	2762	2766	2774	rBV2	84331	145048	1.94%	0.411%
14	19.809	2837	2843	2851	rBV	5870547	7495705	100.00%	21.242%
15	20.492	2956	2959	2966	rVB	126141	154530	2.06%	0.438%
16	21.068	3053	3057	3063	rBV3	150836	234272	3.13%	0.664%
17	21.356	3101	3106	3113	rBV	2036400	2591917	34.58%	7.345%
18	23.632	3487	3493	3508	rVB2	1436224	2840514	37.90%	8.050%

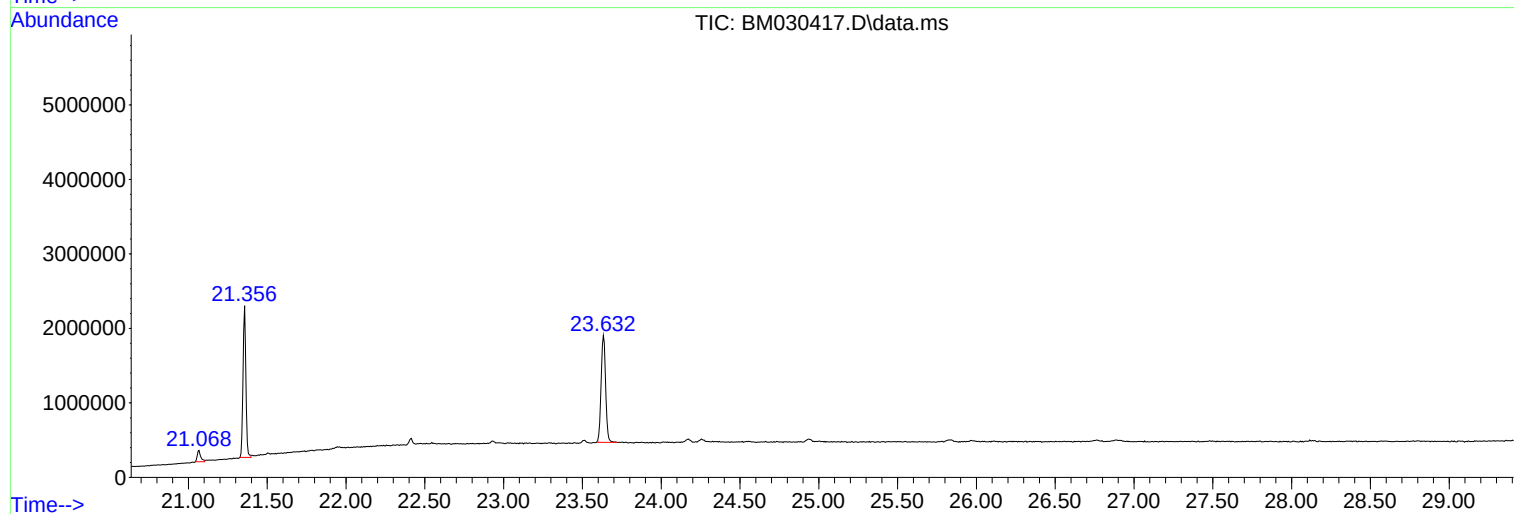
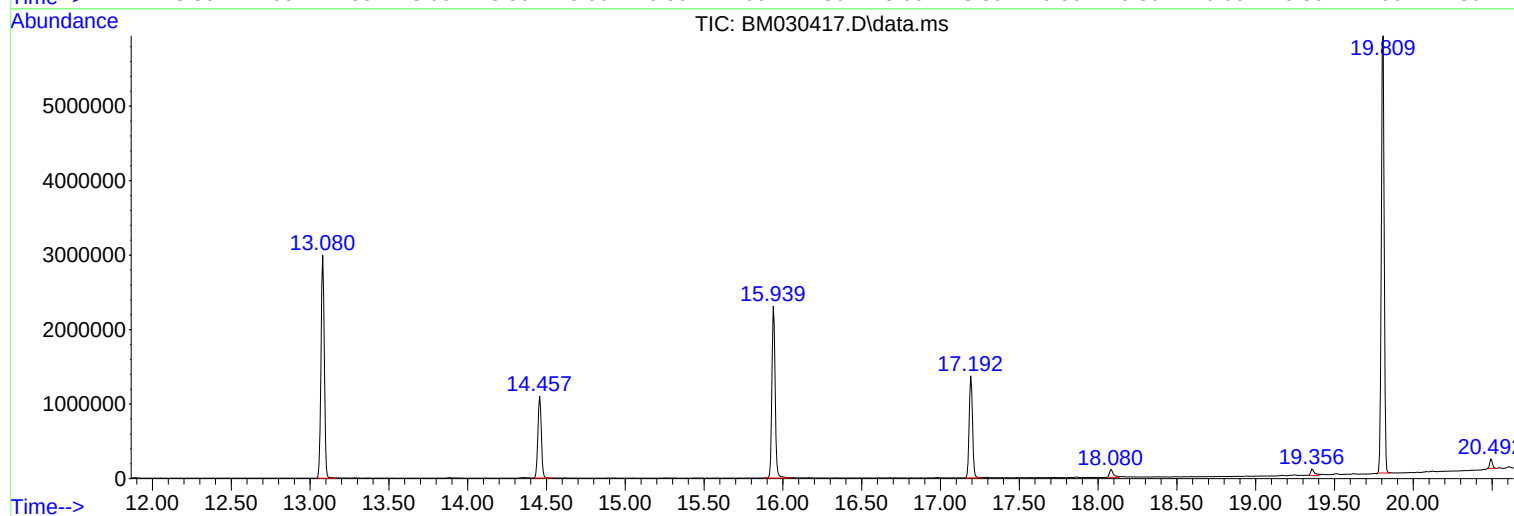
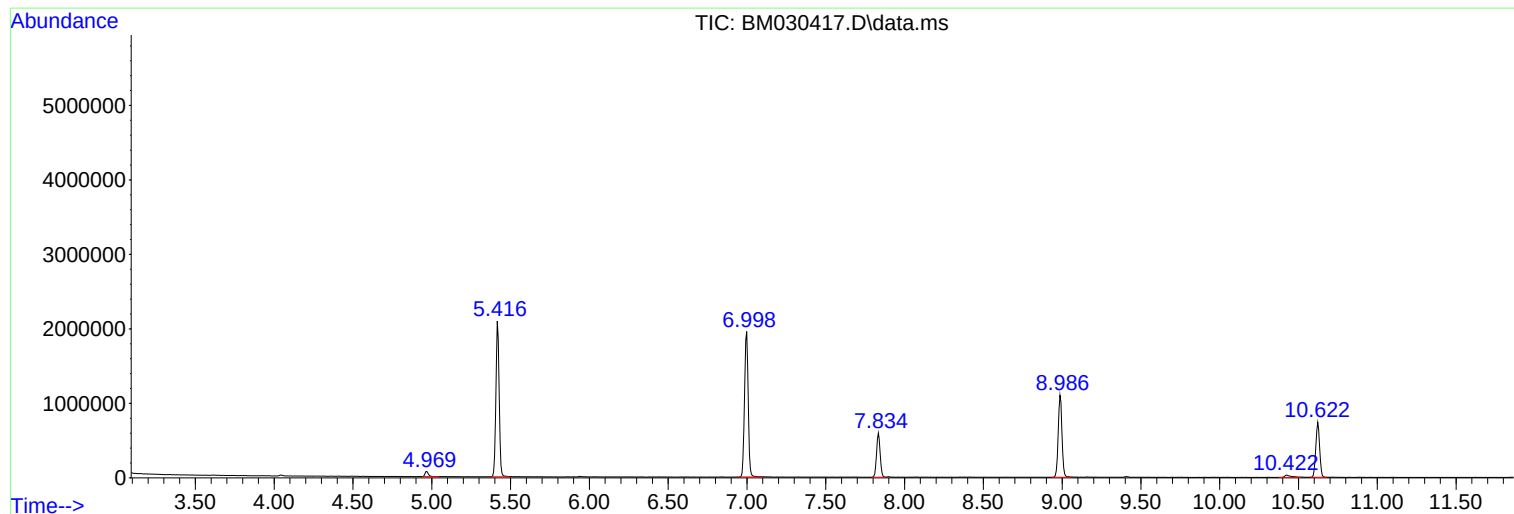
Sum of corrected areas: 35287232

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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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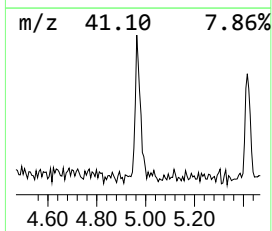
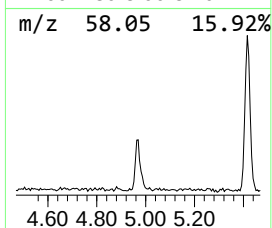
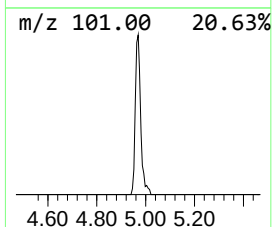
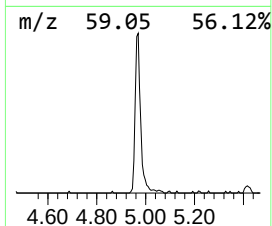
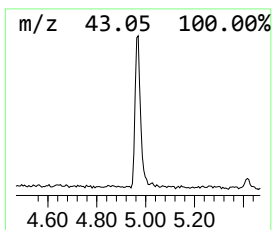
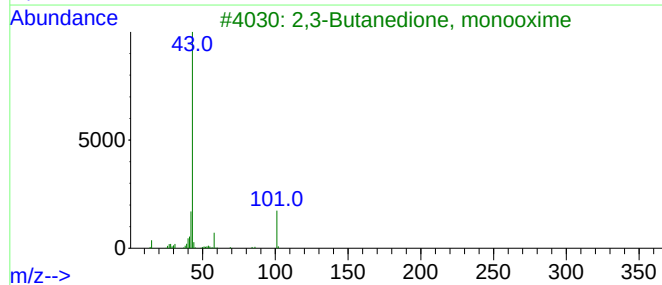
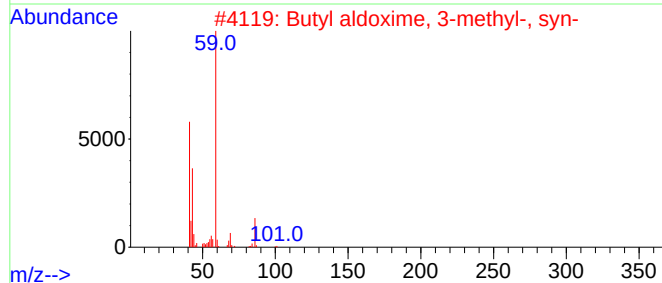
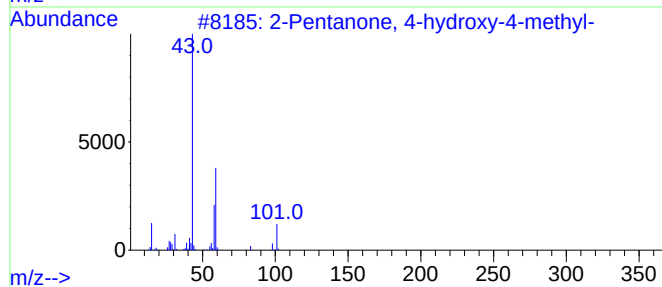
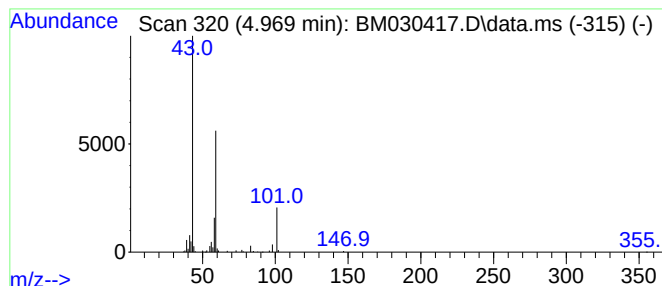
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TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.969	2.34 ng	109472	1,4-Dichlorobenzene-d4	7.833

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50		
2	Butyl aldoxime, 3-methyl-, syn-	101	C5H11NO	005780-40-5	17		
3	2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9		
4	(+)-4-Amino-4,5-dihydro-2(3H)-f...	101	C4H7NO2	016504-58-8	9		
5	5-Hexen-2-one	98	C6H10O	000109-49-9	9		



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-...	4.969	2.3	ng	109472	1	7.833	937067	20.0