

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF060923\
 Data File : BF133691.D
 Acq On : 10 Jun 2023 00:22
 Operator : CG\JU
 Sample : PB153327BL
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB153327BL

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_F\Methods\8270-BF060923.M
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BF133691.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.093	507	511	519	rBV	35363	50953	1.66%	0.281%
2	5.469	569	575	578	rBV	2909932	3035058	98.99%	16.755%
3	6.475	740	746	749	rBV	2622396	3051089	99.52%	16.844%
4	6.840	804	808	811	rBV	651821	497349	16.22%	2.746%
5	7.410	898	905	907	rBV	1785050	1878440	61.27%	10.370%
6	8.122	1022	1026	1029	rBV	792958	591596	19.30%	3.266%
7	9.204	1205	1210	1213	rBV	3271394	3065873	100.00%	16.926%
8	9.881	1321	1325	1328	rBV	626338	540911	17.64%	2.986%
9	10.669	1455	1459	1463	rBV	1761485	1583039	51.63%	8.739%
10	11.369	1574	1578	1582	rBV	611584	504666	16.46%	2.786%
11	12.963	1844	1849	1852	rBV	2985811	2619948	85.46%	14.464%
12	14.016	2024	2028	2032	rBV	409312	360853	11.77%	1.992%
13	15.486	2273	2278	2289	rVB	209621	334086	10.90%	1.844%

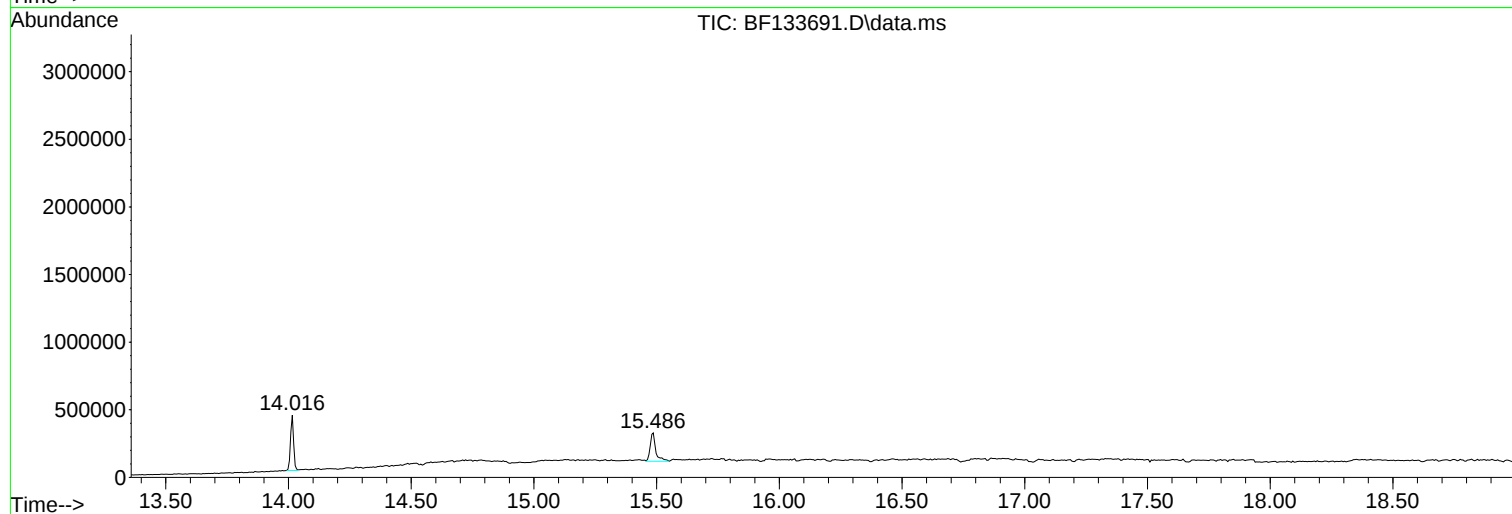
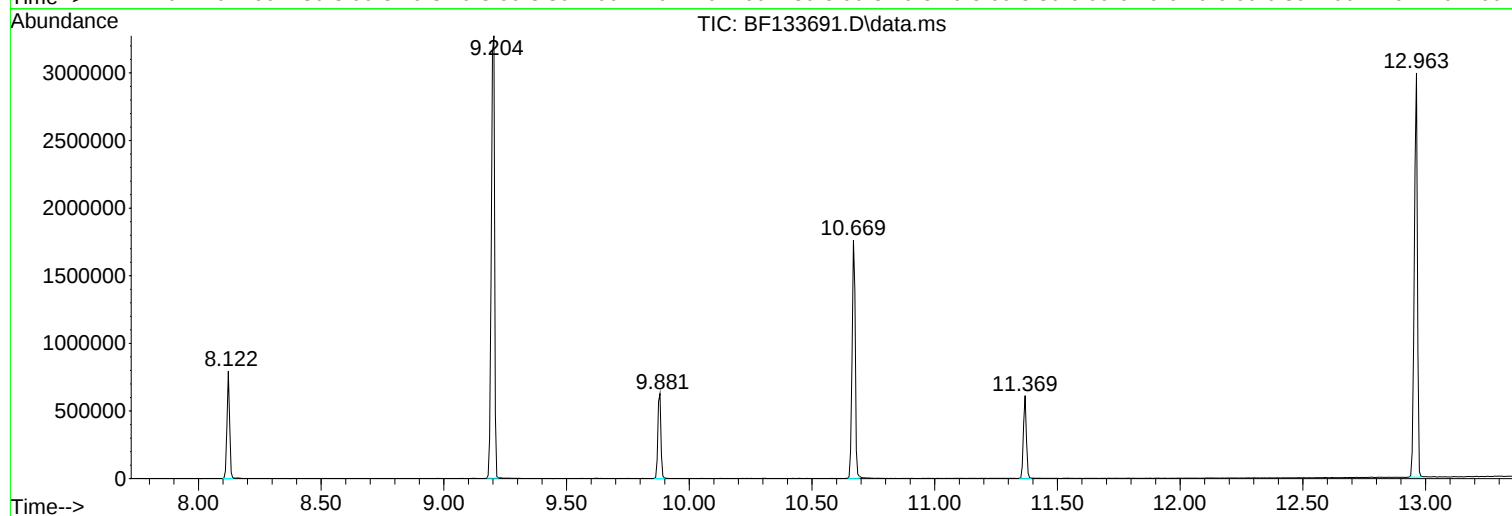
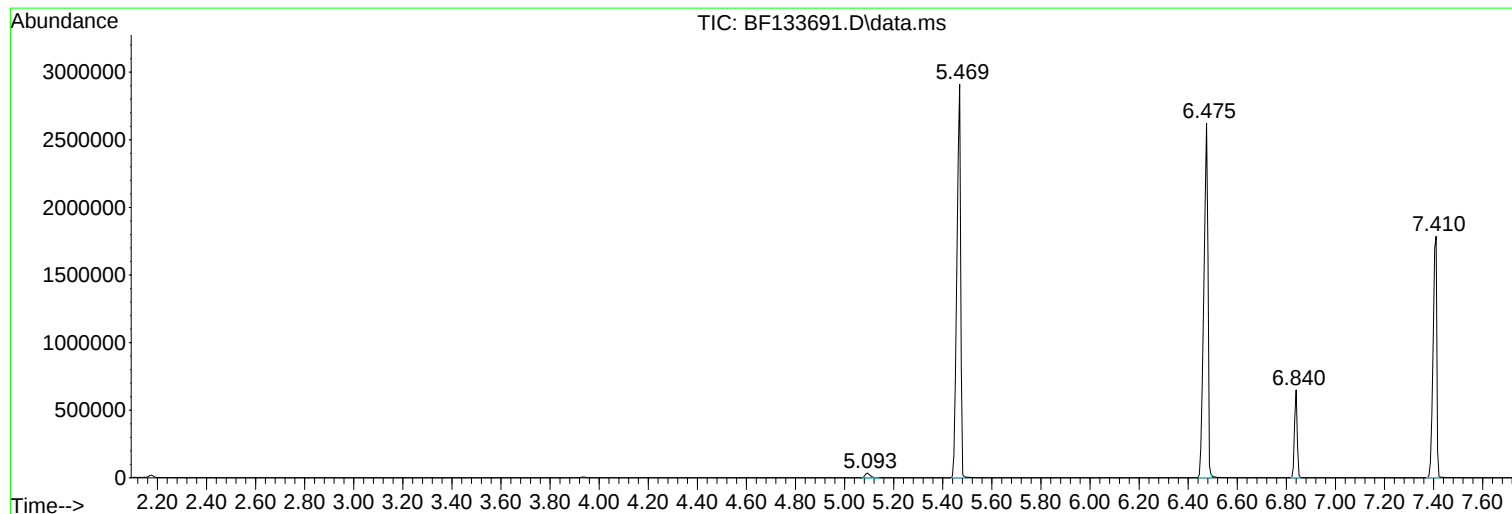
Sum of corrected areas: 18113861

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF060923.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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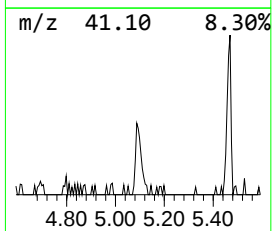
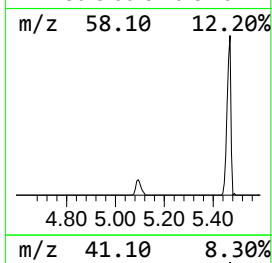
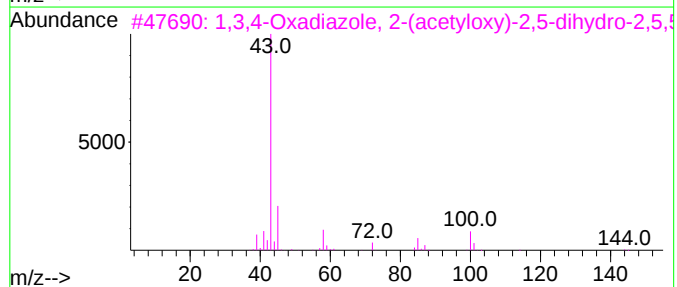
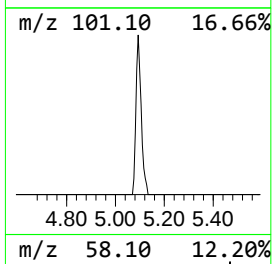
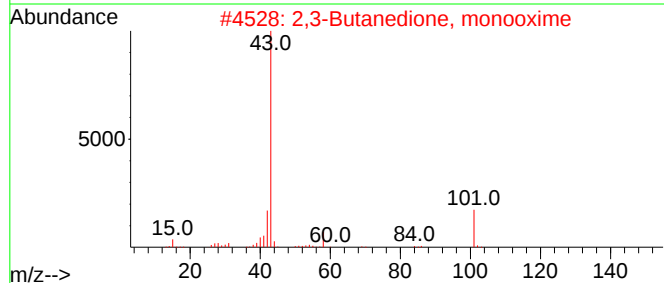
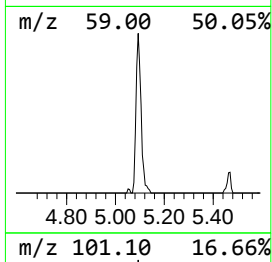
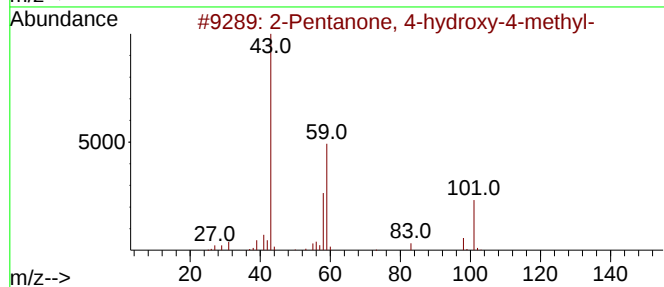
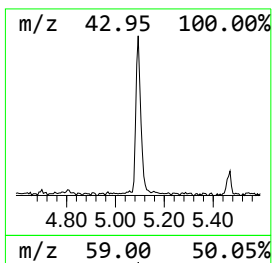
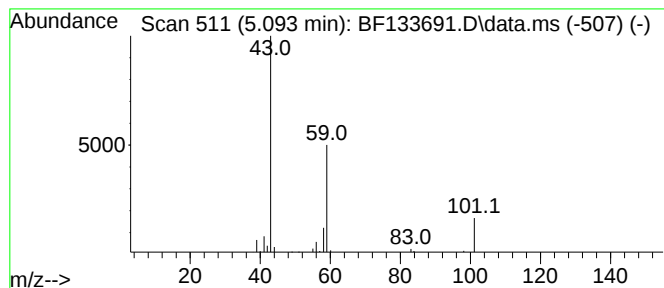
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.093	2.05 ng	50953	1,4-Dichlorobenzene-d4	6.840

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	25	
3	1,3,4-Oxadiazole, 2-(acetyloxy)-...	172	C7H12N2O3	066398-57-0	17	
4	Diacetamide	101	C4H7NO2	000625-77-4	9	
5	1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	9	



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-...	5.093	2.0	ng	50953	1	6.840	497349	20.0