

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF120522\
 Data File : BF131515.D
 Acq On : 05 Dec 2022 23:28
 Operator : CG\JU
 Sample : PB149384BL
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB149384BL

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-BF120522.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BF131515.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.181	521	526	535	rBV	69897	116623	1.50%	0.343%
2	5.545	582	588	591	rBV	3100822	3667894	47.29%	10.787%
3	6.551	752	759	761	rBV	2939456	3620714	46.68%	10.648%
4	6.922	818	822	825	rBV	917082	760532	9.81%	2.237%
5	7.492	912	919	922	rBV	2129307	2370684	30.56%	6.972%
6	8.210	1035	1041	1044	rBV	1074926	978322	12.61%	2.877%
7	9.292	1218	1225	1228	rBV	5122578	5447381	70.23%	16.020%
8	9.975	1334	1341	1344	rBV	1479569	1243572	16.03%	3.657%
9	10.769	1469	1476	1479	rBV	3892255	3904311	50.34%	11.482%
10	11.469	1589	1595	1598	rBV	1907136	1614879	20.82%	4.749%
11	13.068	1859	1867	1870	rBV	7199993	7756259	100.00%	22.811%
12	14.127	2042	2047	2050	rBV	1467166	1381881	17.82%	4.064%
13	15.651	2299	2306	2319	rBV	832073	1139797	14.70%	3.352%

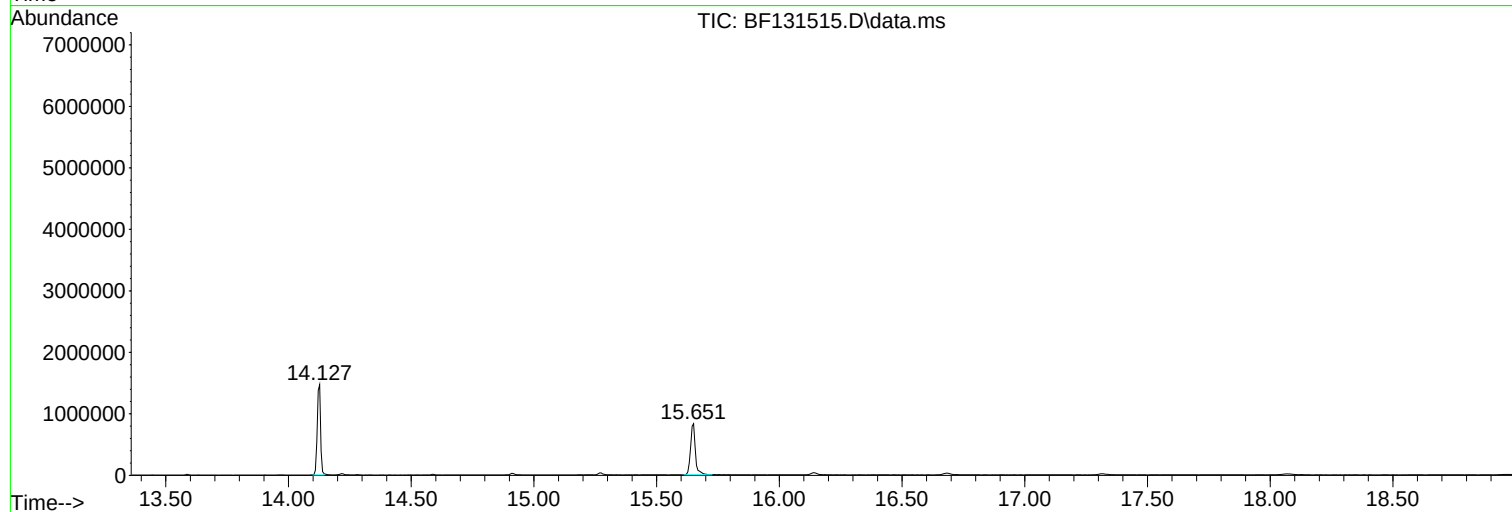
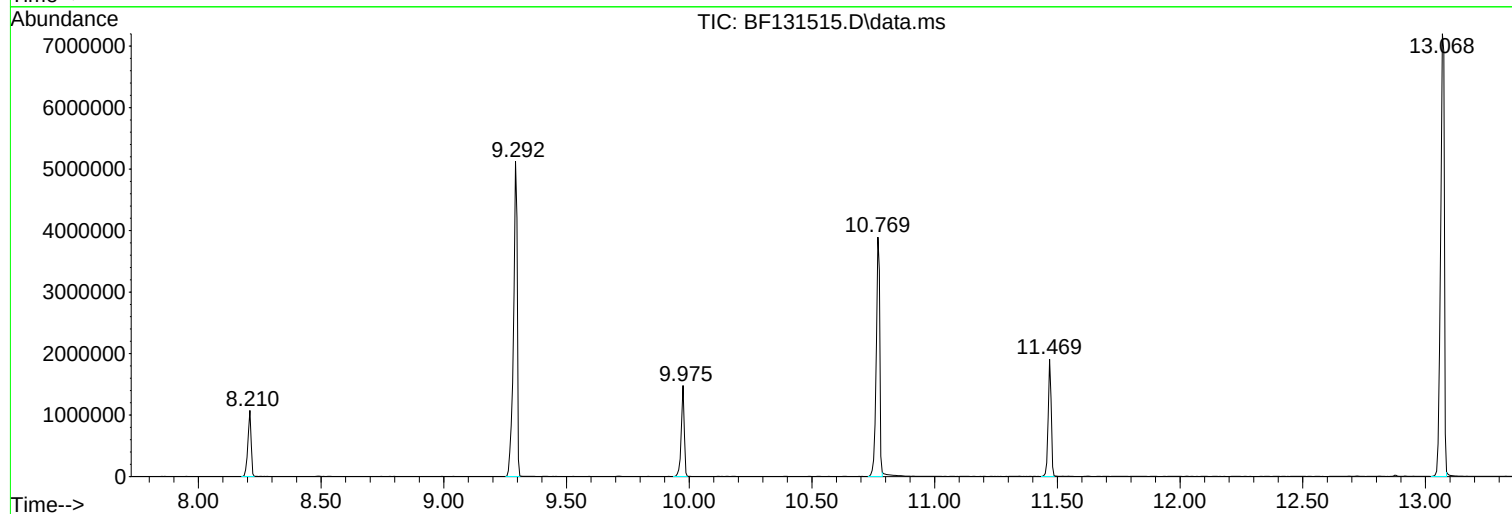
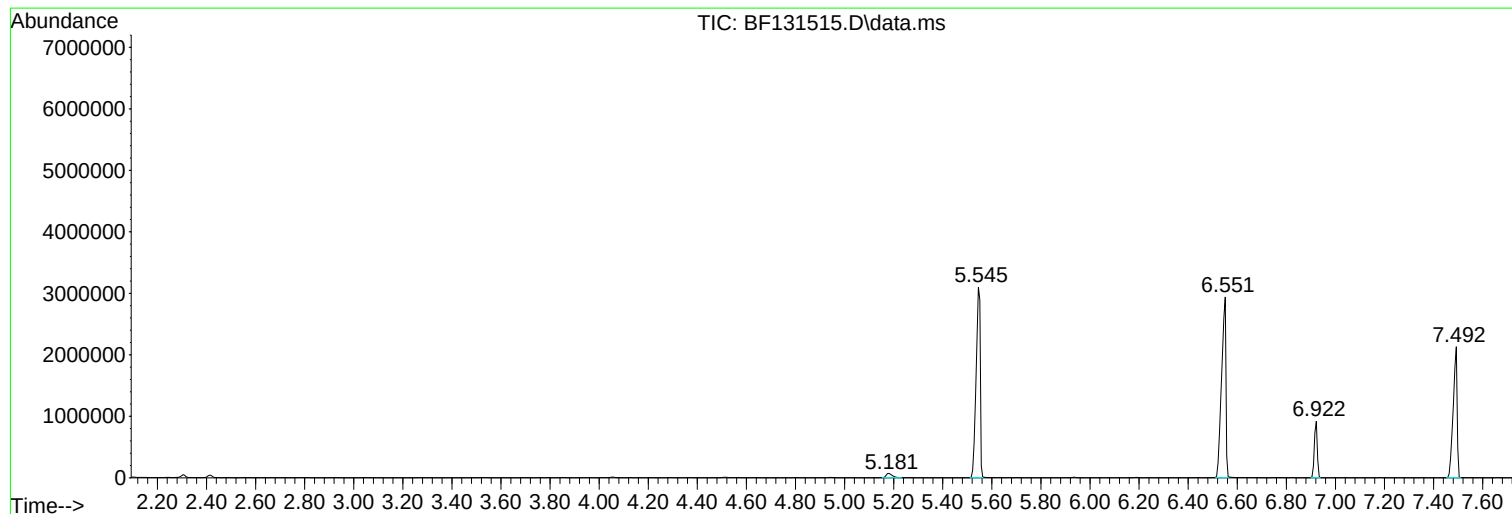
Sum of corrected areas: 34002849

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF120522.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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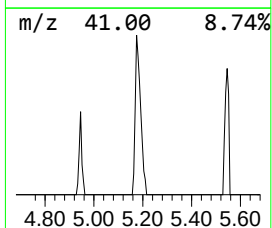
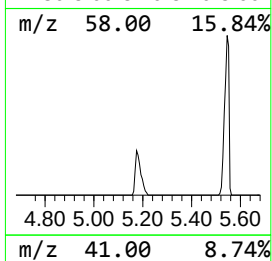
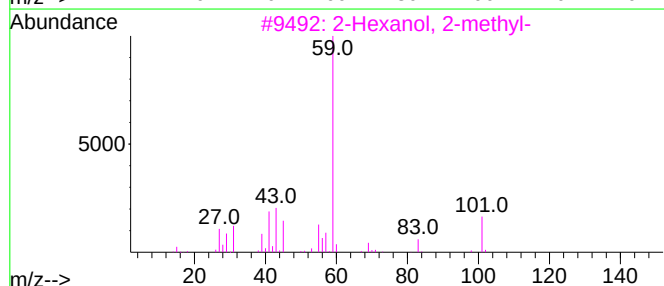
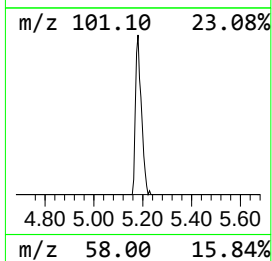
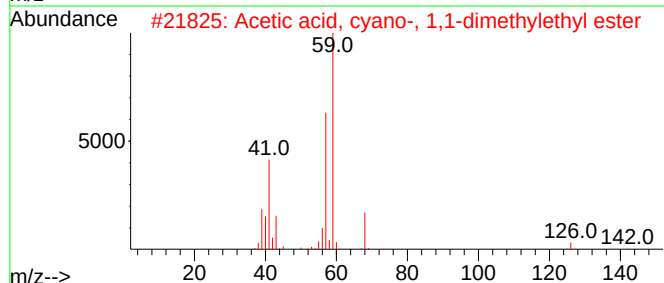
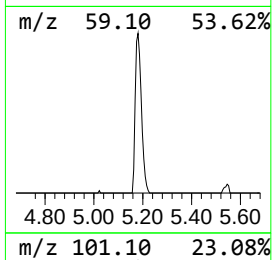
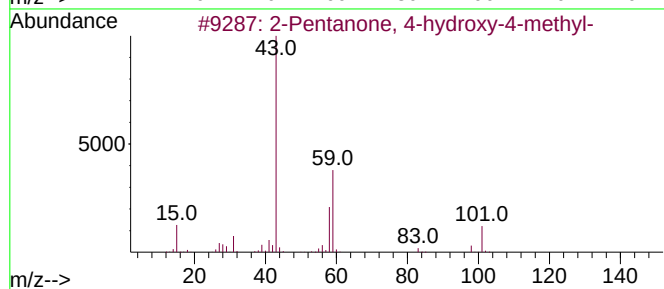
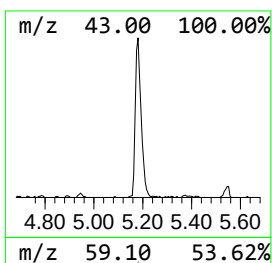
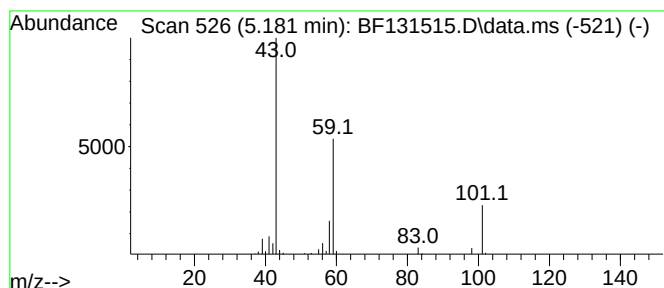
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.181	3.07 ng	116623	1,4-Dichlorobenzene-d4	6.922

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50		
2	Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	25		
3	2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	25		
4	2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	17		
5	Butyl aldoxime, 3-methyl-, syn-	101	C5H11NO	005780-40-5	12		



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-...	5.181	3.1	ng	116623	1	6.922	760532	20.0