

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF112222\
 Data File : BF131322.D
 Acq On : 22 Nov 2022 10:17
 Operator : CG\JU
 Sample : PB148922BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB148922BL

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-BF112122.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BF131322.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.352	31	45	50	rBV	118684	174113	3.67%	0.634%
2	5.204	525	530	546	rVB	338150	491065	10.35%	1.788%
3	5.587	588	595	598	rBV	2942999	3462358	73.01%	12.609%
4	6.581	757	764	767	rBV	2835199	3435508	72.44%	12.511%
5	6.963	825	829	832	rBV	931011	729524	15.38%	2.657%
6	7.528	919	925	928	rBV	1965950	2297082	48.44%	8.365%
7	8.251	1043	1048	1052	rBV	1042770	952680	20.09%	3.469%
8	9.333	1226	1232	1235	rBV	4143355	4257408	89.77%	15.504%
9	10.016	1342	1348	1352	rBV	1366048	1162279	24.51%	4.233%
10	10.810	1476	1483	1487	rBV	2680679	2767294	58.35%	10.078%
11	11.516	1599	1603	1606	rVB	1425654	1259316	26.55%	4.586%
12	13.116	1869	1875	1878	rBV	4739280	4742319	100.00%	17.270%
13	14.168	2050	2054	2058	rBV	1165725	1007953	21.25%	3.671%
14	15.710	2310	2316	2321	rBV	591489	721161	15.21%	2.626%

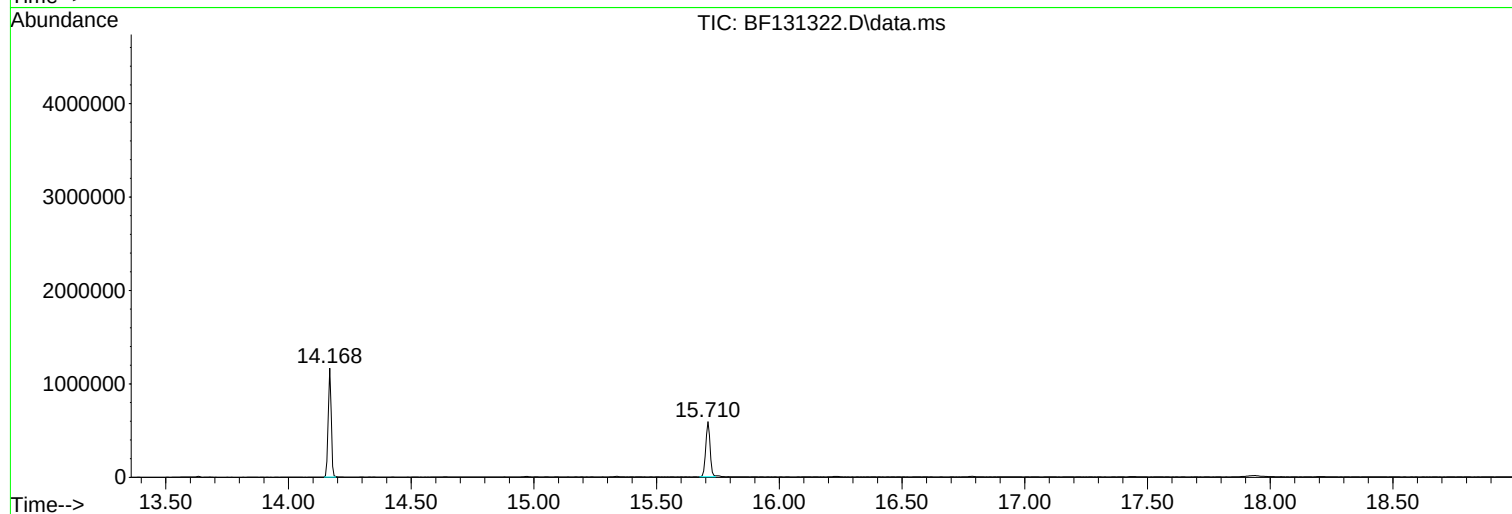
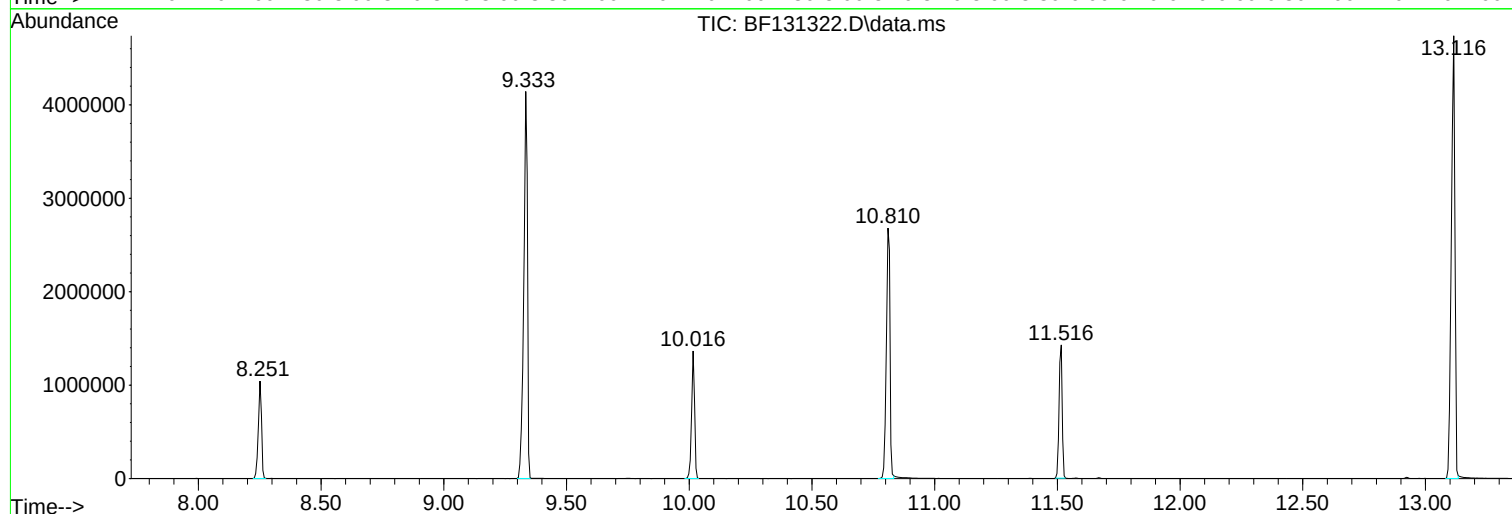
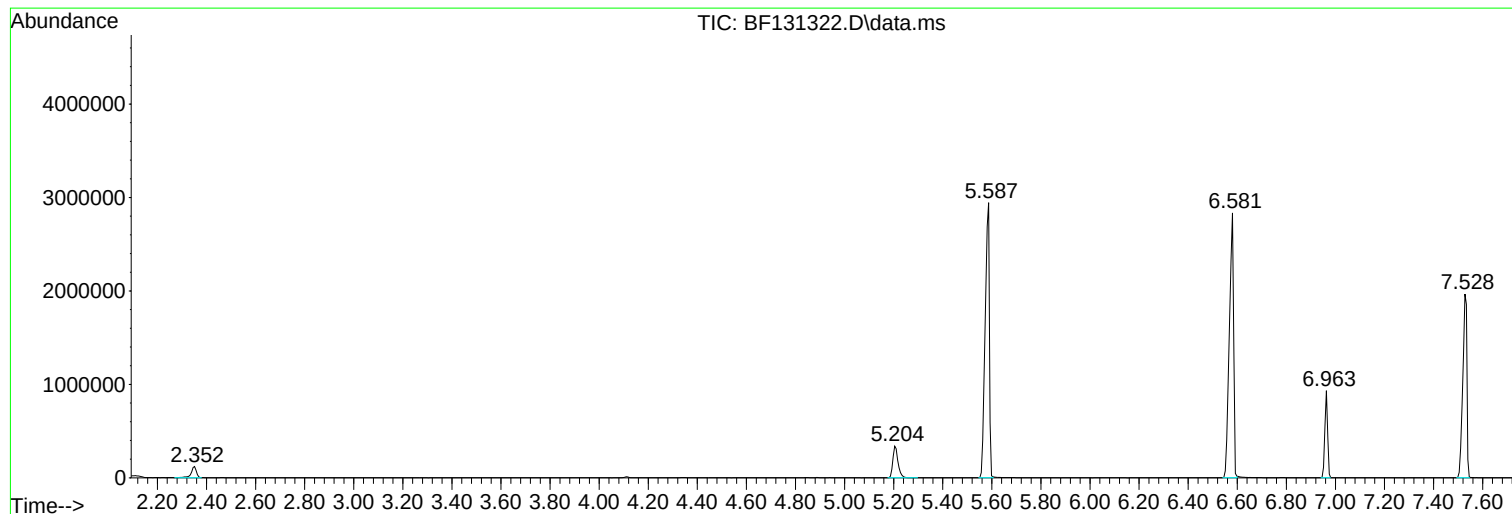
Sum of corrected areas: 27460060

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF112122.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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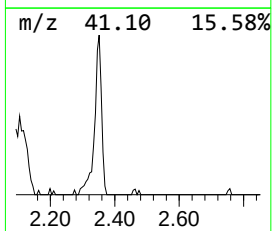
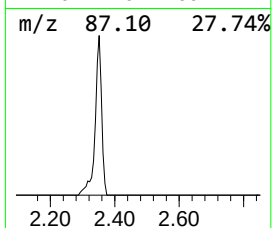
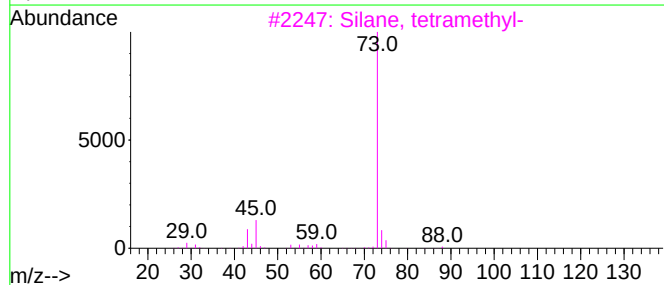
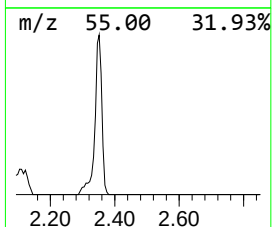
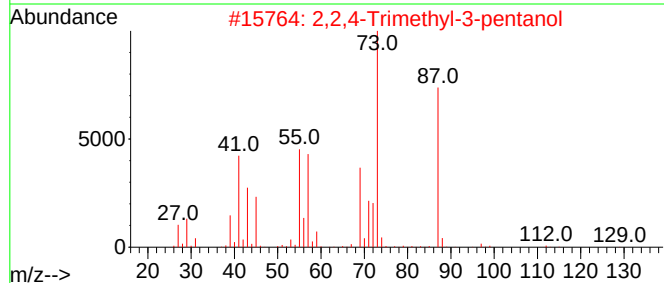
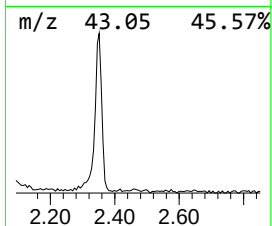
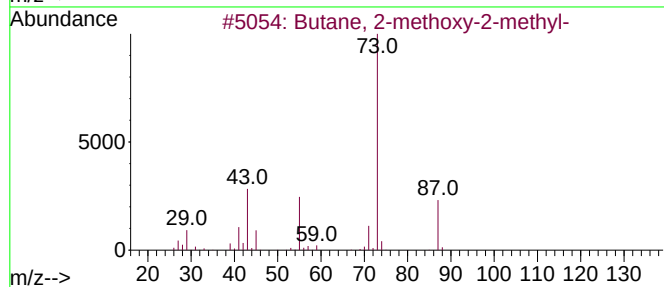
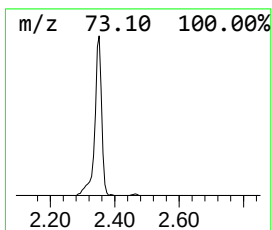
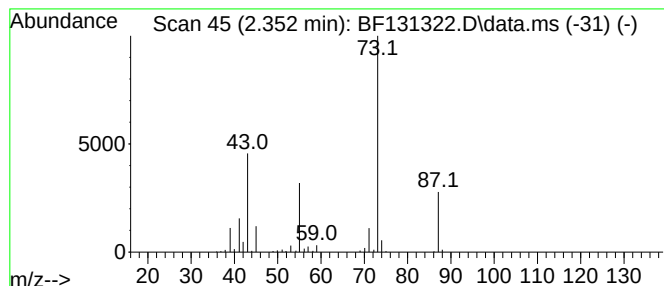
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TIC Integration Parameters: LSCINT.P

Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.352	4.77 ng	174113	1,4-Dichlorobenzene-d4	6.963

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	78
2			2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	40
3			Silane, tetramethyl-	88	C4H12Si	000075-76-3	38
4			1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	9
5			Silane, ethyltrimethyl-	102	C5H14Si	003439-38-1	9



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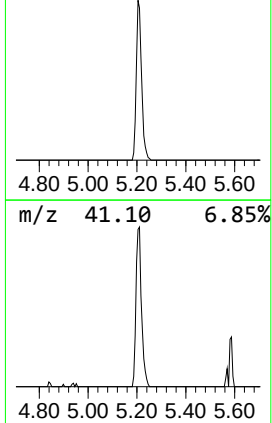
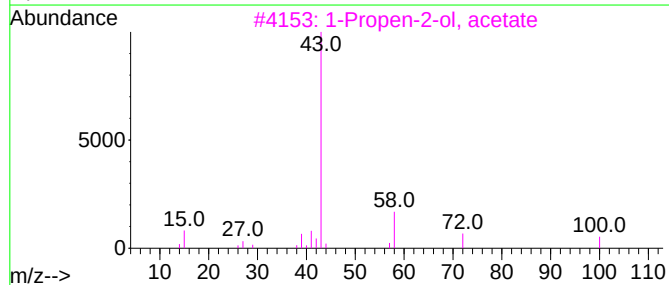
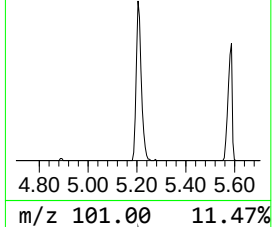
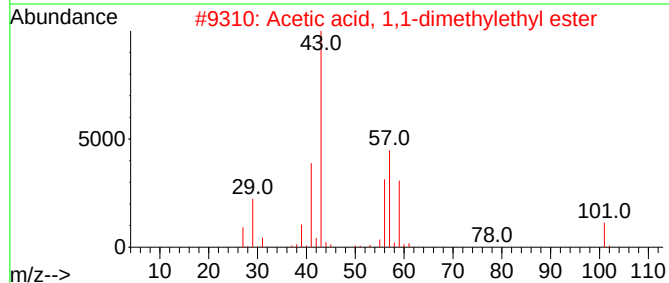
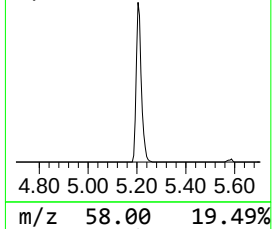
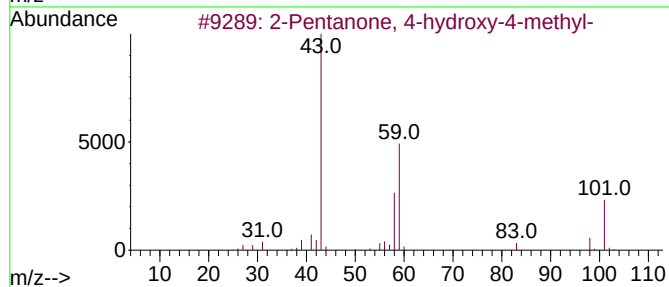
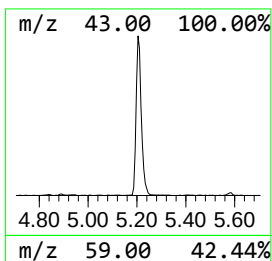
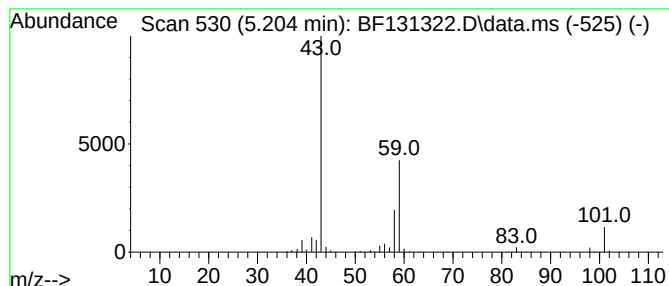
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Peak Number 2 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.204	13.46 ng	491065	1,4-Dichlorobenzene-d4	6.963

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	83		
2	Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	36		
3	1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	16		
4	Acetone	58	C3H6O	000067-64-1	16		
5	Diacetamide	101	C4H7NO2	000625-77-4	9		



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Butane, 2-metho...	2.352	4.8	ng	174113	1	6.963	729524	20.0
2-Pentanone, 4-...	5.204	13.5	ng	491065	1	6.963	729524	20.0