

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF032521\
 Data File : BF123462.D
 Acq On : 25 Mar 2021 15:23
 Operator : JU/CG
 Sample : PB135267BL
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB135267BL

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-BF032321.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BF123462.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.281	23	33	38	rVV	105172	143441	1.77%	0.274%
2	2.387	40	51	57	rVB	289524	376021	4.63%	0.718%
3	5.146	516	520	528	rVB	261806	343166	4.23%	0.655%
4	5.528	579	585	588	rBV	6586519	6817388	83.98%	13.020%
5	6.522	748	754	757	rBV	5863799	7006376	86.31%	13.381%
6	6.898	814	818	821	rBV	1927117	1538447	18.95%	2.938%
7	7.463	907	914	916	rBV	4102086	4600433	56.67%	8.786%
8	8.181	1031	1036	1040	rBV	2233405	2025594	24.95%	3.869%
9	9.263	1213	1220	1223	rBV	7712414	7643134	94.15%	14.597%
10	9.939	1330	1335	1339	rBV	2597469	2306746	28.42%	4.405%
11	10.733	1463	1470	1473	rBV	5344216	5090945	62.71%	9.723%
12	11.428	1582	1588	1592	rBV	2533305	2337911	28.80%	4.465%
13	13.027	1853	1860	1863	rBV	8246818	8117866	100.00%	15.504%
14	14.080	2034	2039	2042	rBV	2155280	2132075	26.26%	4.072%
15	15.574	2285	2293	2298	rBV	1495723	1881103	23.17%	3.593%

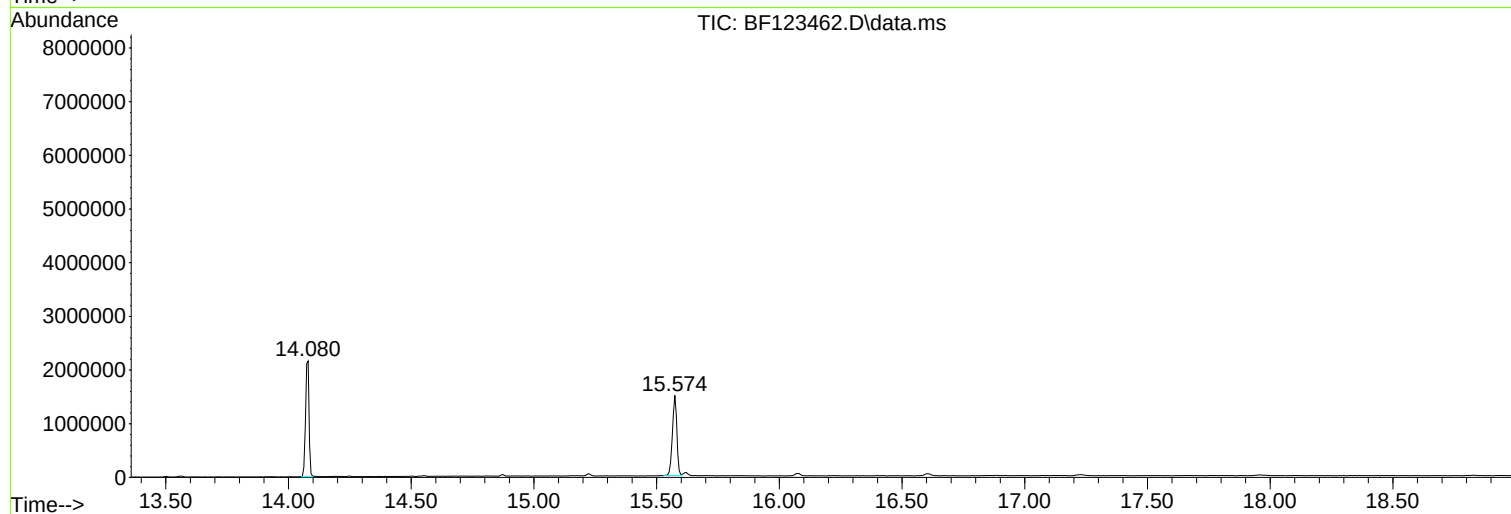
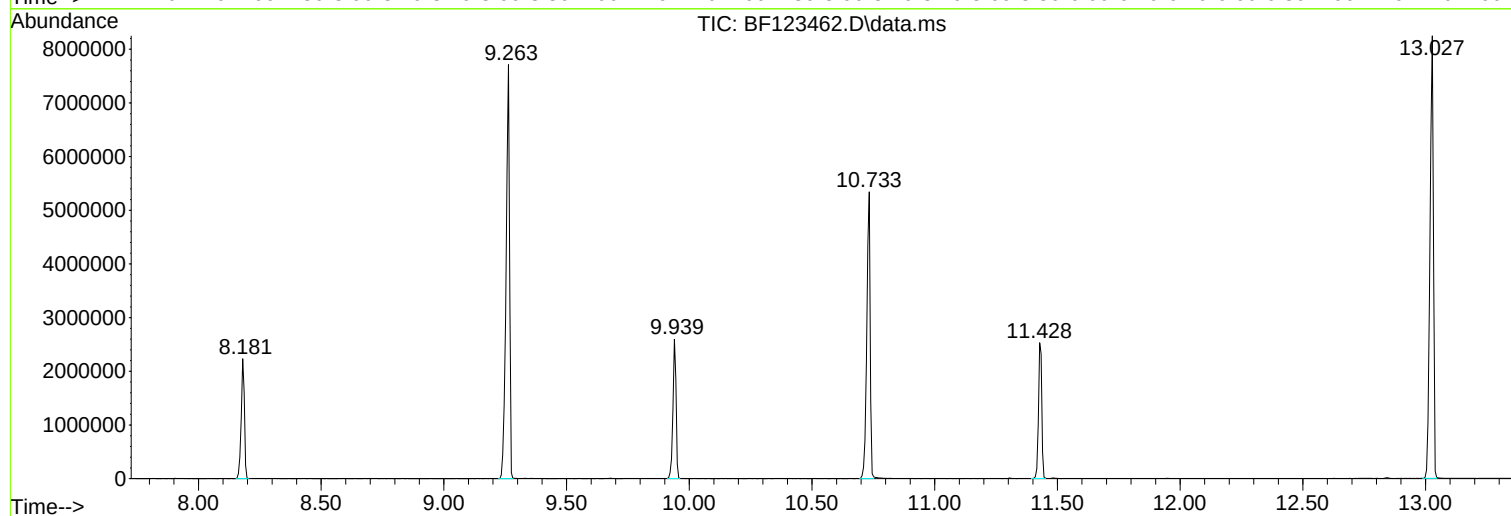
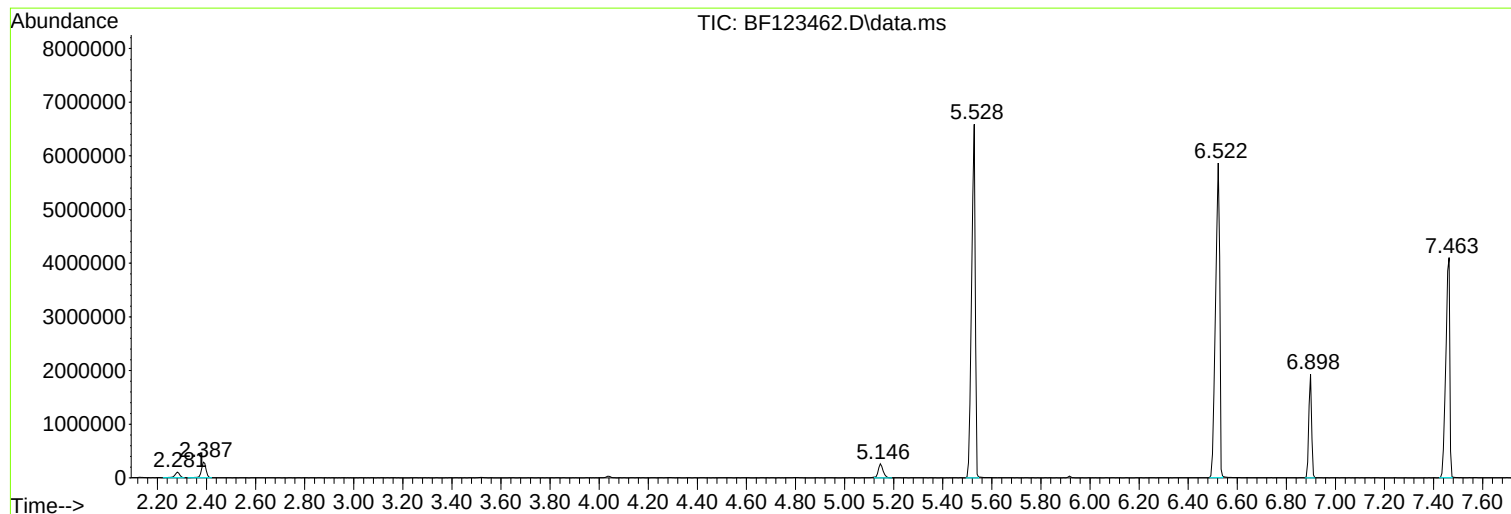
Sum of corrected areas: 52360646

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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF032321.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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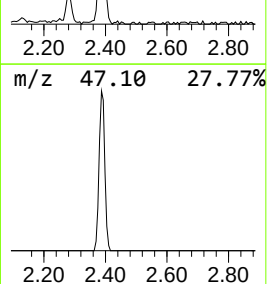
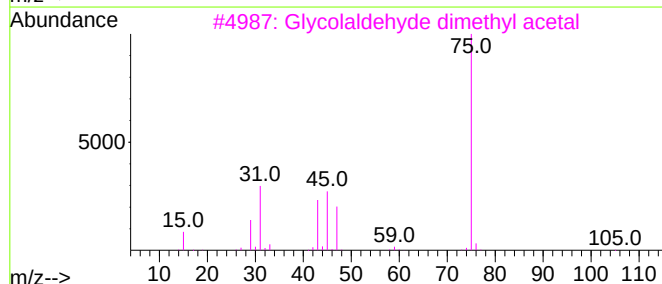
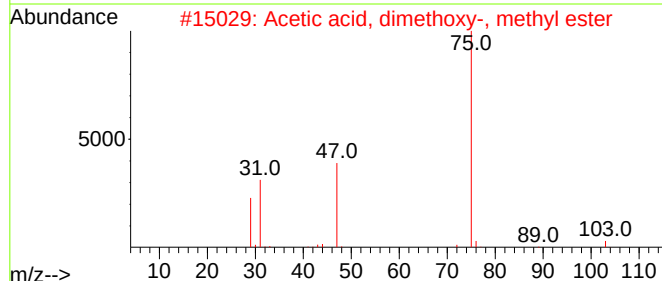
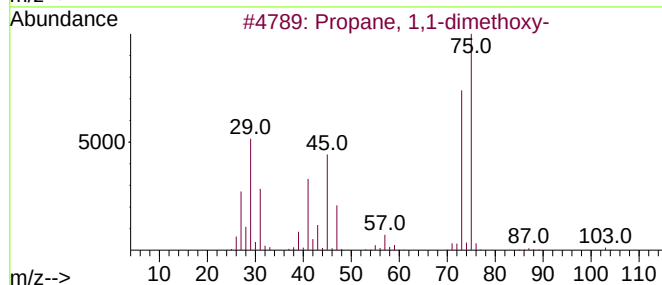
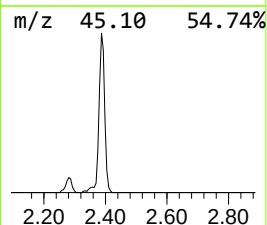
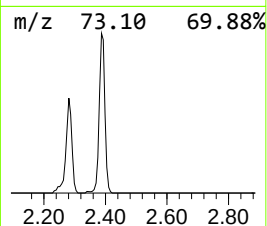
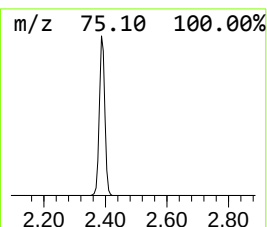
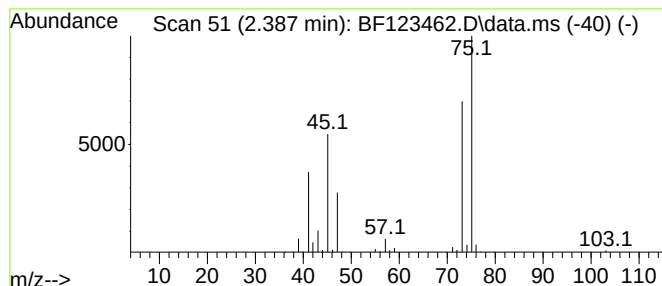
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Peak Number 1 Propane, 1,1-dimethoxy- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.387	4.89 ng	376021	1,4-Dichlorobenzene-d4	6.898

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Propane, 1,1-dimethoxy-	104	C5H12O2	004744-10-9	83
2			Acetic acid, dimethoxy-, methyl ...	134	C5H10O4	000089-91-8	50
3			Glycolaldehyde dimethyl acetal	106	C4H10O3	030934-97-5	45
4			Aminoacetaldehyde dimethyl acetal	105	C4H11NO2	022483-09-6	36
5			Ethane, 1,1,2-trimethoxy-	120	C5H12O3	024332-20-5	36



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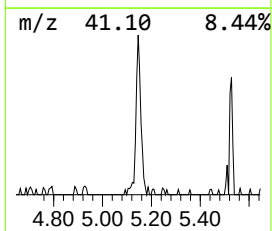
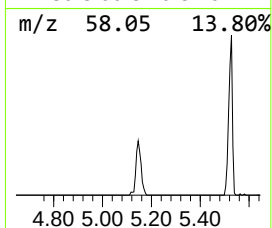
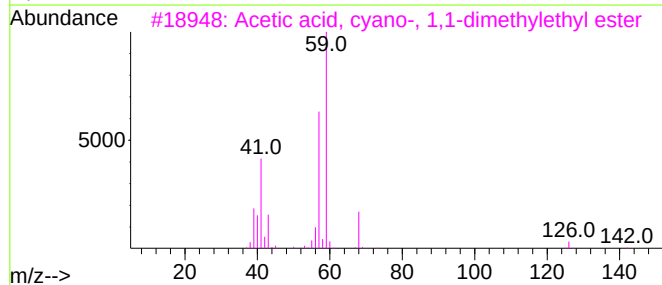
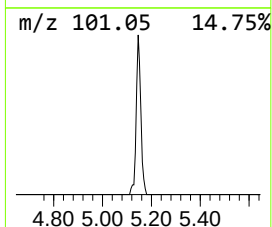
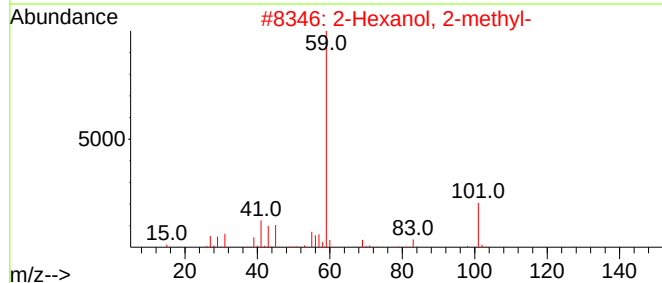
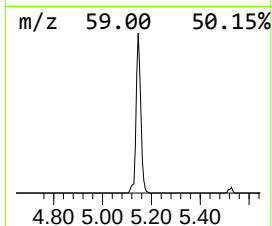
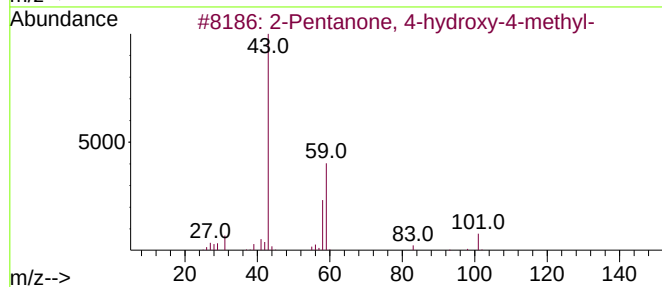
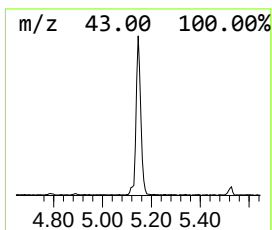
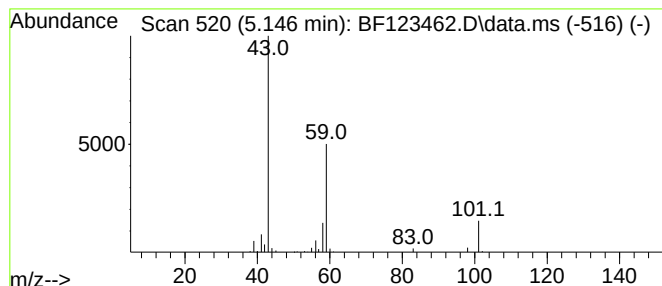
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Peak Number 2 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.146	4.46 ng	343166	1,4-Dichlorobenzene-d4	6.898

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50
2			2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	33
3			Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	25
4			3-Hexanol, 4-ethyl-	130	C8H18O	019780-44-0	9
5			3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Propane, 1,1-di...	2.387	4.9	ng	376021	1	6.898	1538450	20.0
2-Pentanone, 4-...	5.146	4.5	ng	343166	1	6.898	1538450	20.0