

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF040120\
 Data File : BF119687.D
 Acq On : 2 Apr 2020 00:49
 Operator : MA/SJ
 Sample : PB127927BL
 Misc :
 ALS Vial : 34 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB127927BL

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-BF031820.M
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.240	22	26	35	rVB	126205	164750	1.81%	0.280%
2	5.045	498	503	514	rBV	226880	333064	3.65%	0.566%
3	5.434	563	569	572	rBV	6579115	6996545	76.74%	11.889%
4	6.440	734	740	743	rBV	5520844	6900150	75.68%	11.725%
5	6.810	799	803	806	rBV	1674928	1319204	14.47%	2.242%
6	6.969	825	830	833	rBV	6024262	5559332	60.98%	9.447%
7	7.375	893	899	902	rBV	4298936	4601938	50.47%	7.820%
8	8.092	1016	1021	1025	rBV	1821634	1737399	19.06%	2.952%
9	9.175	1199	1205	1208	rBV	7685542	8340897	91.48%	14.173%
10	9.851	1314	1320	1324	rBV	2307788	2030065	22.27%	3.450%
11	10.645	1448	1455	1458	rBV	4933254	5402963	59.26%	9.181%
12	11.339	1567	1573	1576	rBV	2603238	2247484	24.65%	3.819%
13	12.933	1836	1844	1847	rBV	7907895	9117309	100.00%	15.492%
14	13.980	2018	2022	2025	rBV	2316302	2040990	22.39%	3.468%
15	15.439	2263	2270	2274	rBV	1598382	2058560	22.58%	3.498%

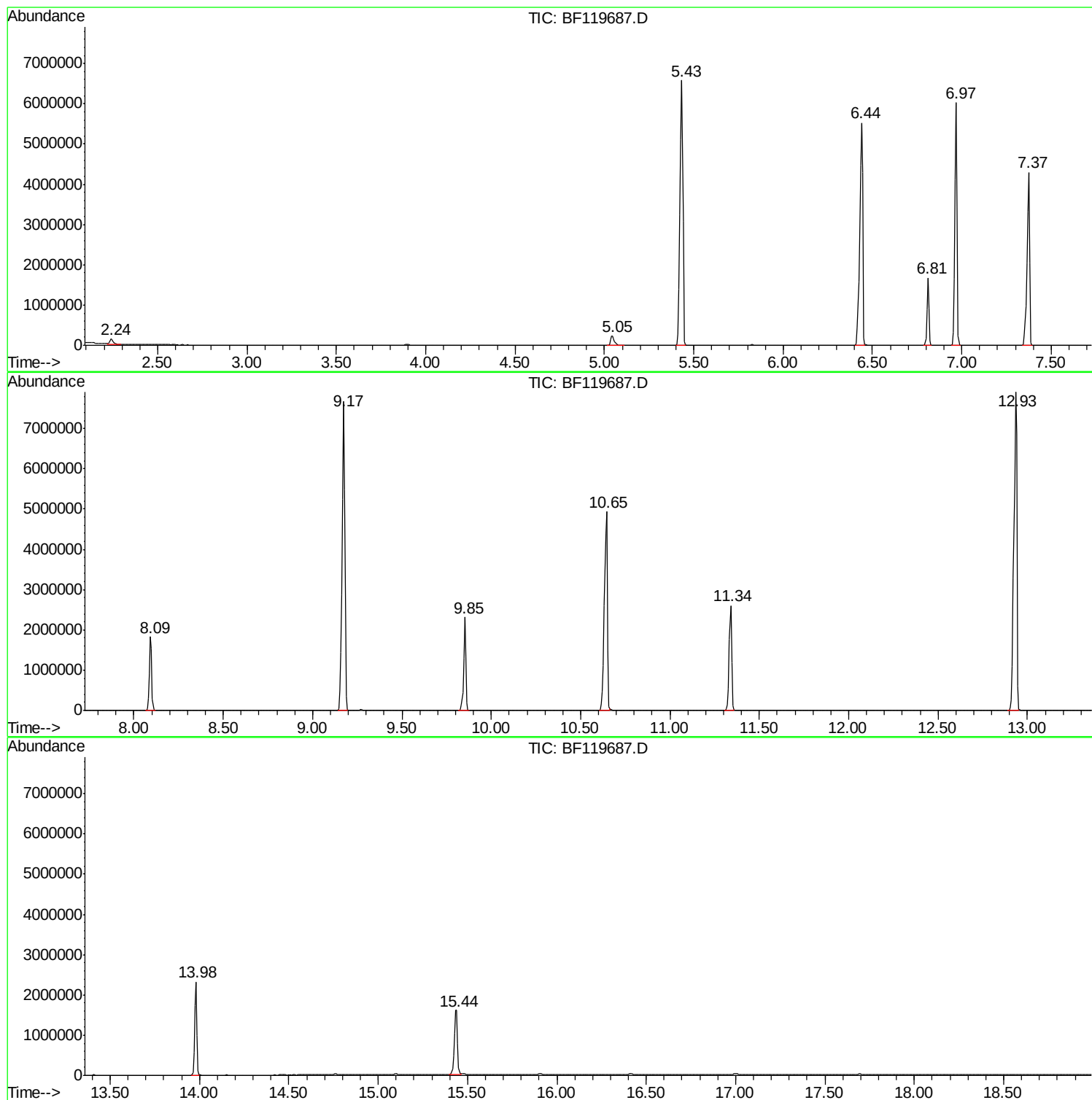
Sum of corrected areas: 58850650

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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P



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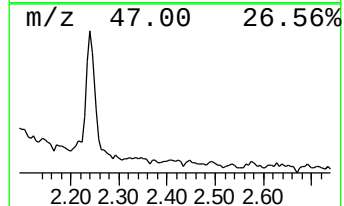
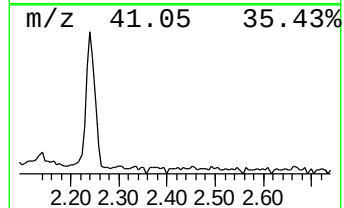
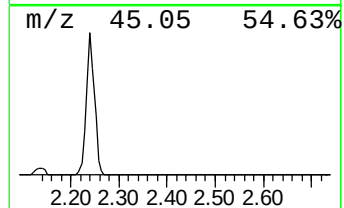
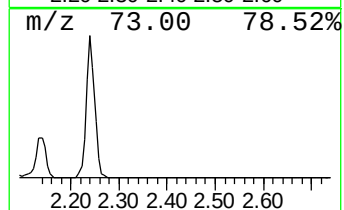
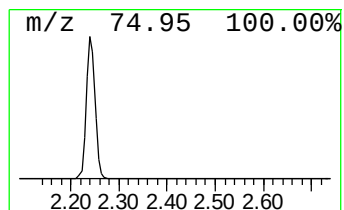
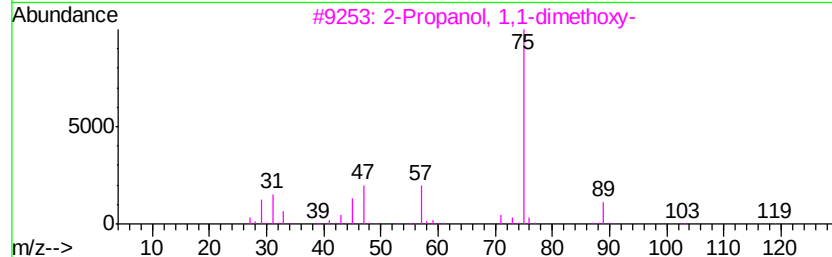
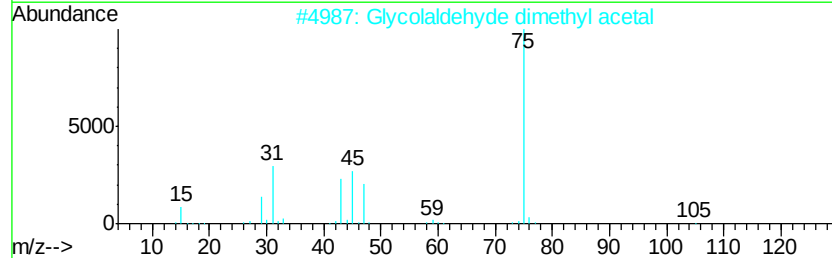
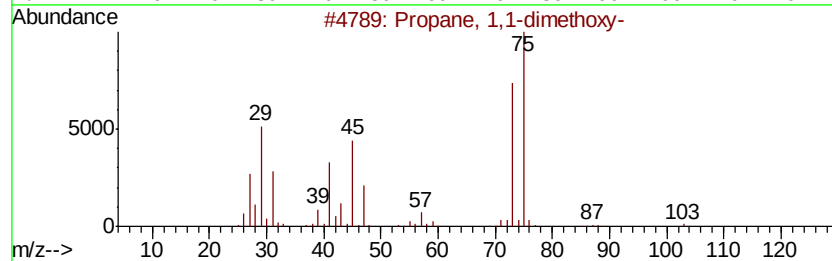
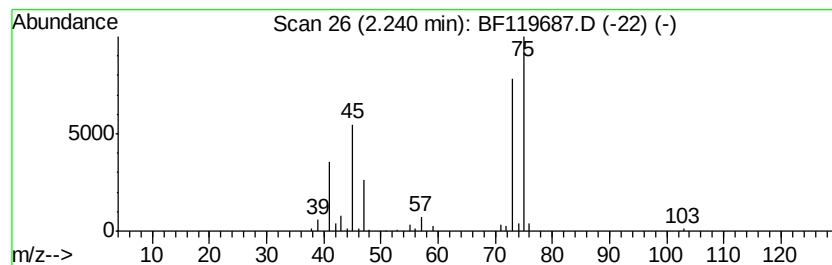
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TIC Integration Parameters: LSCINT.P

Peak Number 1 Propane, 1,1-dimethoxy- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.24	2.50 ng	164750	1,4-Dichlorobenzene-d4	6.81

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Propane, 1,1-dimethoxy-	104	C5H12O2	004744-10-9	83
2		Glycolaldehyde dimethyl acetal	106	C4H10O3	030934-97-5	42
3		2-Propanol, 1,1-dimethoxy-	120	C5H12O3	042919-42-6	40
4		Acetic acid, dimethoxy-, methyl ...	134	C5H10O4	000089-91-8	38
5		Aminoacetaldehyde dimethyl acetal	105	C4H11NO2	022483-09-6	28



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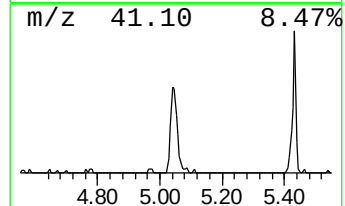
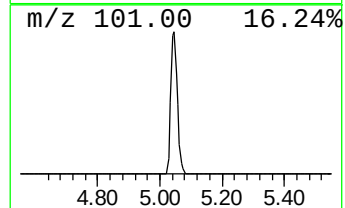
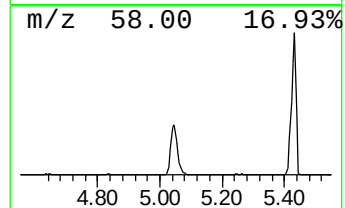
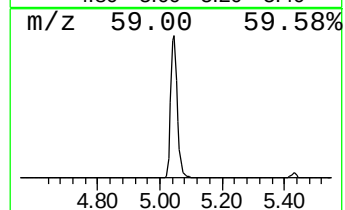
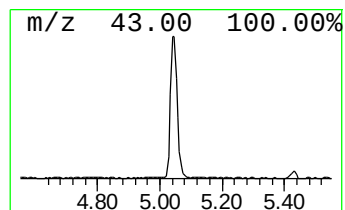
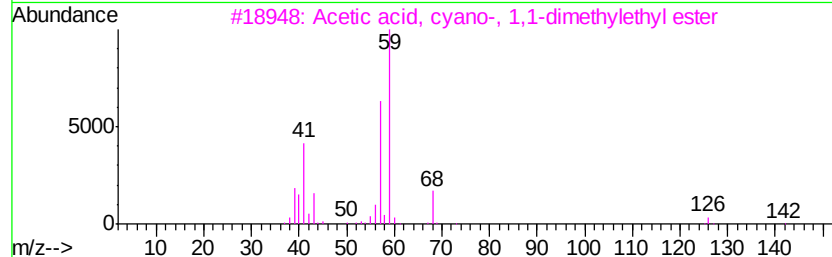
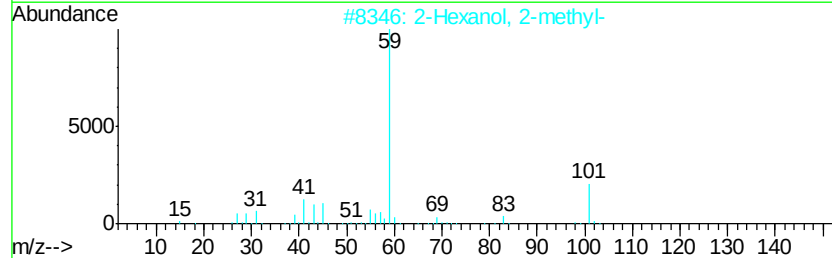
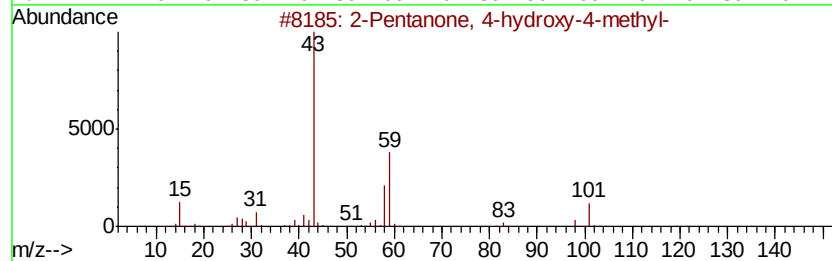
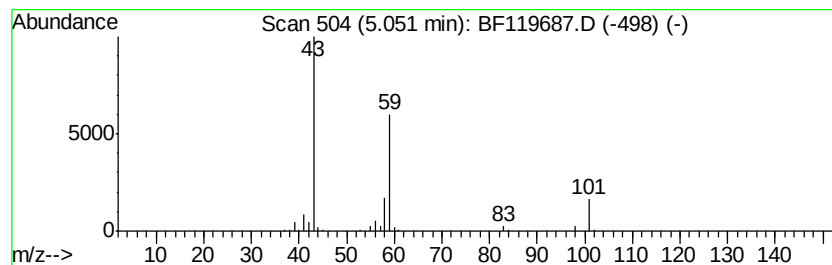
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Peak Number 2 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.05	5.05 ng	333064	1,4-Dichlorobenzene-d4	6.81

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	59
2			2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	33
3			Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	32
4			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	16
5			Pentane, 2-methoxy-	102	C6H14O	006795-88-6	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Propane, 1,1-dime...	2.24	2.5	ng	164750	1	6.81	1319200	20.0
2-Pentanone, 4-hy...	5.05	5.0	ng	333064	1	6.81	1319200	20.0