

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF032720\
 Data File : BF119575.D
 Acq On : 27 Mar 2020 18:55
 Operator : CG/JU
 Sample : PB127867BL
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB127867BL

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	2.269	24	31	36	rBV	178749	227062	2.30%	0.331%
2	5.069	499	507	515	rBV	161488	217427	2.20%	0.317%
3	5.451	566	572	575	rBV	8267133	8971443	90.89%	13.084%
4	6.457	737	743	746	rBV	7275616	8678856	87.93%	12.657%
5	6.833	803	807	810	rBV	1871816	1555453	15.76%	2.268%
6	6.992	829	834	837	rBV	7030804	6809544	68.99%	9.931%
7	7.398	896	903	905	rBV	4938783	5865038	59.42%	8.554%
8	8.116	1020	1025	1028	rBV	2295563	2039037	20.66%	2.974%
9	9.198	1202	1209	1212	rBV	9648344	9822828	99.52%	14.326%
10	9.874	1318	1324	1327	rBV	2852232	2446198	24.78%	3.568%
11	10.663	1451	1458	1462	rBV	6112085	6487670	65.73%	9.462%
12	11.363	1571	1577	1580	rBV	2684083	2307080	23.37%	3.365%
13	12.957	1842	1848	1851	rBV	9654991	9870131	100.00%	14.395%
14	14.004	2021	2026	2029	rBV	2002995	1830340	18.54%	2.669%
15	15.462	2268	2274	2280	rBV	1137565	1439924	14.59%	2.100%

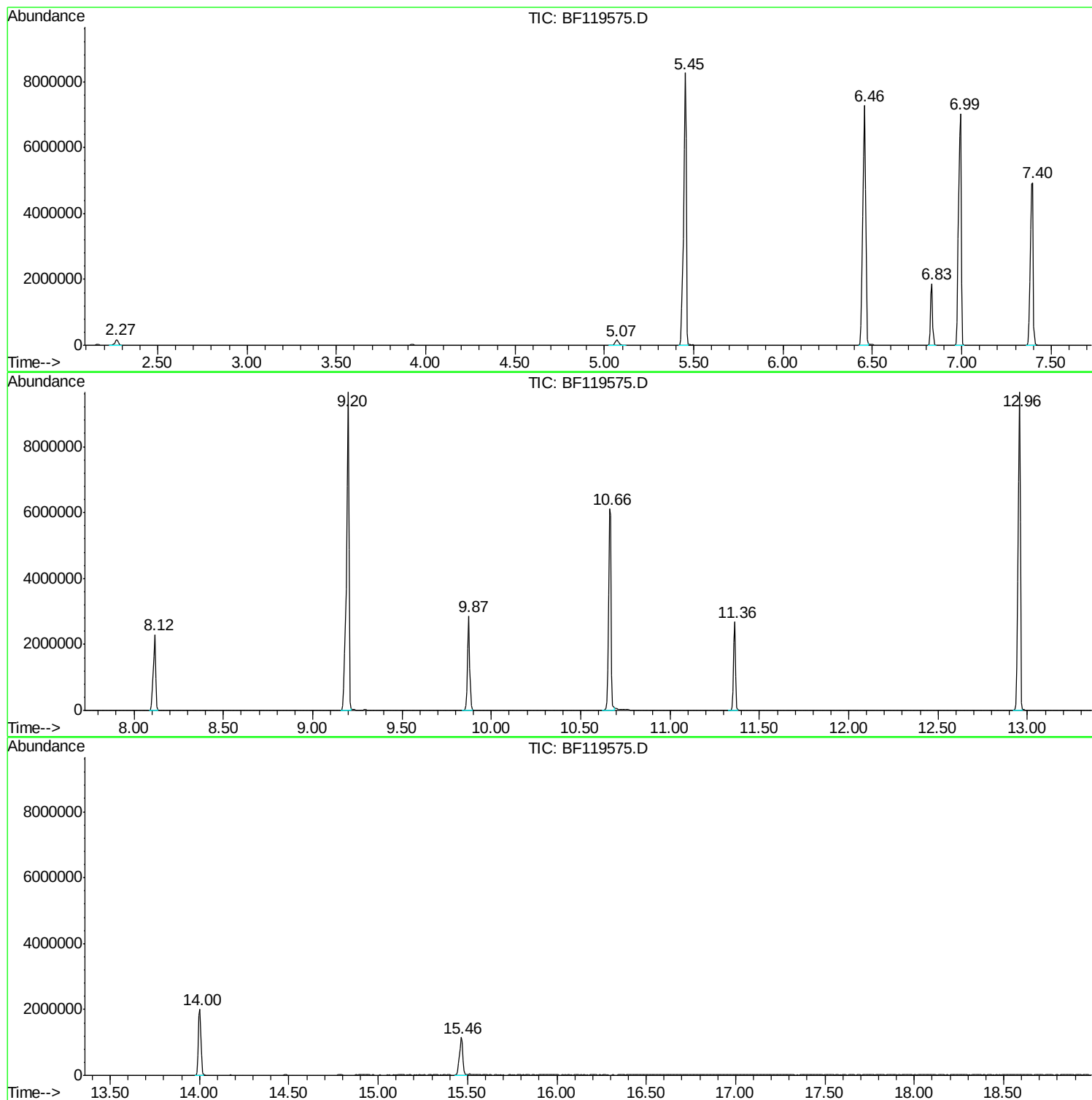
Sum of corrected areas: 68568031

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



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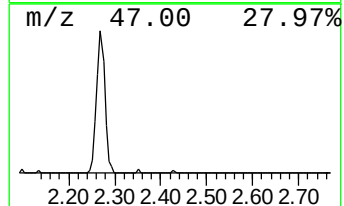
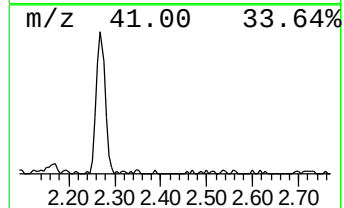
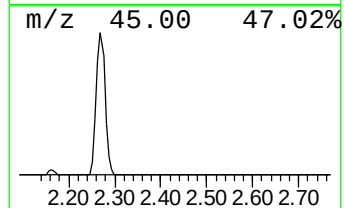
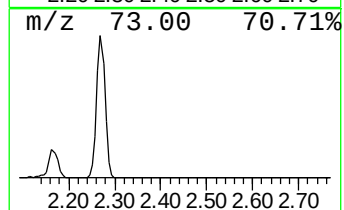
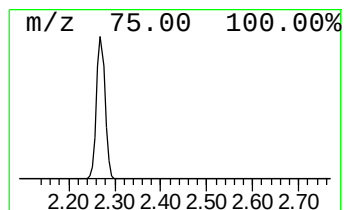
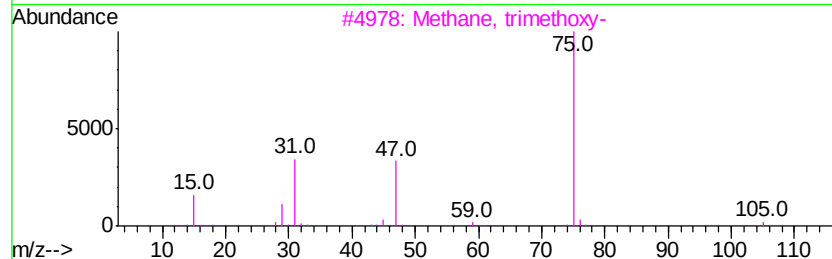
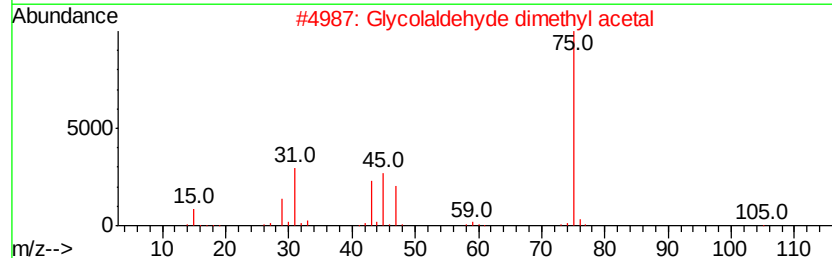
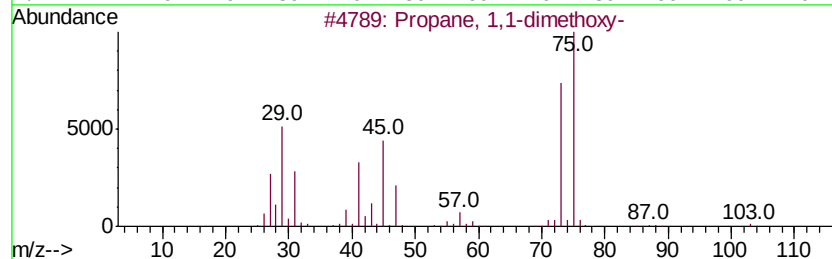
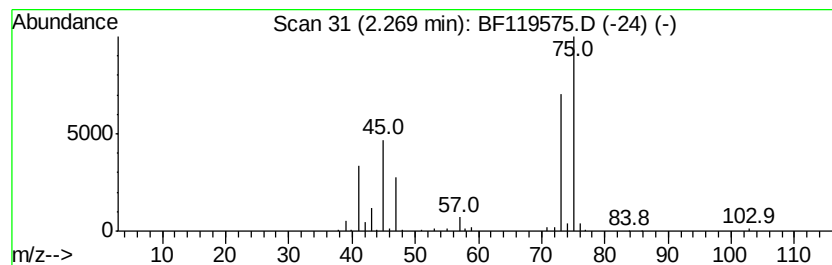
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Peak Number 1 Propane, 1,1-dimethoxy- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.27	2.92 ng	227062	1,4-Dichlorobenzene-d4	6.83

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Propane, 1,1-dimethoxy-	104	C5H12O2	004744-10-9	78
2		Glycolaldehyde dimethyl acetal	106	C4H10O3	030934-97-5	45
3		Methane, trimethoxy-	106	C4H10O3	000149-73-5	40
4		Aminoacetaldehyde dimethyl acetal	105	C4H11NO2	022483-09-6	36
5		Acetic acid, dimethoxy-, methyl ...	134	C5H10O4	000089-91-8	33



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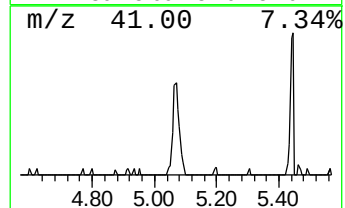
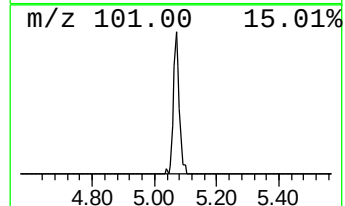
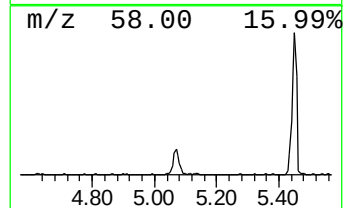
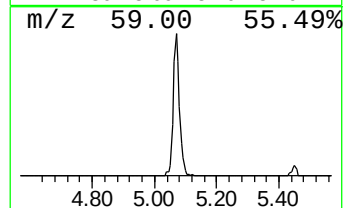
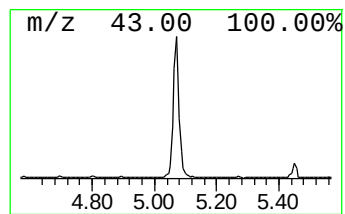
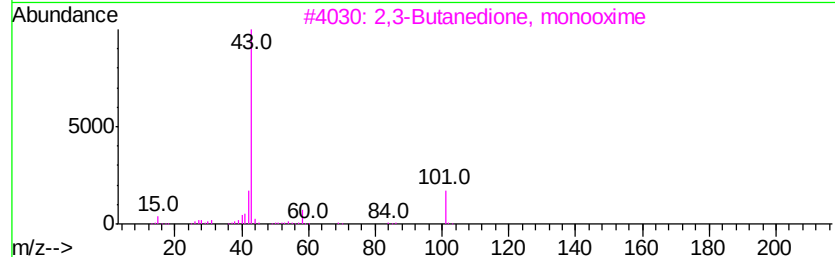
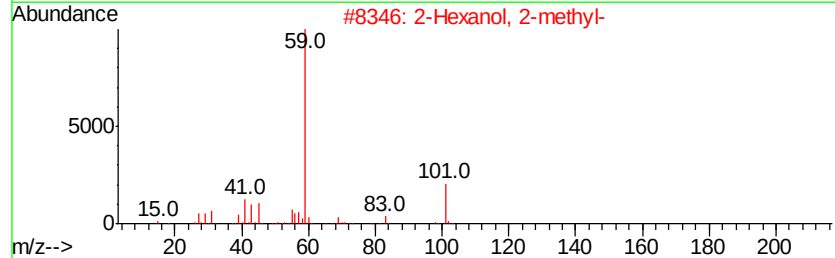
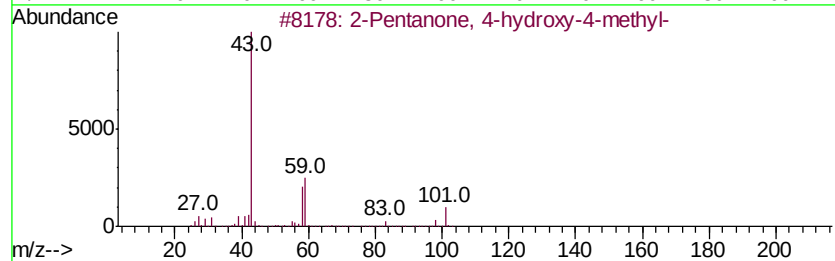
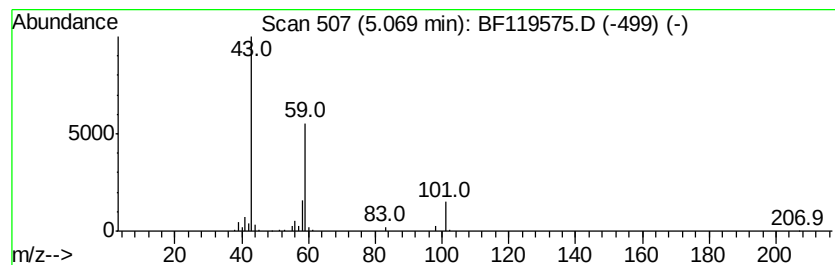
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Peak Number 2 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.07	2.80 ng	217427	1,4-Dichlorobenzene-d4	6.83

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	59
2		2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	28
3		2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	10
4		Propanoic acid, 3-nitro-, methyl...	133	C4H7NO4	020497-95-4	9
5		2-Nonanone, 9-hydroxy-	158	C9H18O2	025368-56-3	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Propane, 1,1-dime...	2.27	2.9	ng	227062	1	6.83	1555450	20.0
2-Pentanone, 4-hy...	5.07	2.8	ng	217427	1	6.83	1555450	20.0