

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF100920\
 Data File : BF121902.D
 Acq On : 9 Oct 2020 12:47
 Operator : JU/CG
 Sample : PB132237BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB132237BL

Integration Parameters: rteint.p
 Integrator: RTE
 Smoothing : ON
 Sampling : 1
 Start Thrs: 0.2
 Stop Thrs : 0
 Filtering: 5
 Min Area: 3 % of largest Peak
 Max Peaks: 100
 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-BF091020.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	2.187	13	17	29	rVB	70565	103842	1.91%	0.337%
2	5.004	492	496	508	rBV	55660	92321	1.70%	0.299%
3	5.392	557	562	565	rBV	3575225	3704049	68.26%	12.015%
4	6.404	729	734	737	rBV	3263622	3722460	68.60%	12.075%
5	6.757	790	794	807	rVB	945293	849282	15.65%	2.755%
6	7.328	885	891	907	rBV	2332455	2839423	52.33%	9.210%
7	8.039	1007	1012	1033	rBV	1083035	1121203	20.66%	3.637%
8	9.116	1189	1195	1207	rBV	4033207	4815889	88.75%	15.621%
9	9.792	1304	1310	1325	rBV	1259942	1286404	23.71%	4.173%
10	10.586	1439	1445	1473	rBV	2853376	3022382	55.70%	9.804%
11	11.280	1557	1563	1580	rBV	1235492	1339613	24.69%	4.345%
12	12.868	1826	1833	1836	rBV	4781871	5426116	100.00%	17.601%
13	13.921	2007	2012	2025	rBV	1146780	1260092	23.22%	4.087%
14	15.357	2251	2256	2269	rBV	804364	1131066	20.84%	3.669%
15	15.780	2323	2328	2335	rVB	37904	55876	1.03%	0.181%
16	16.256	2404	2409	2416	rVB	34792	58589	1.08%	0.190%

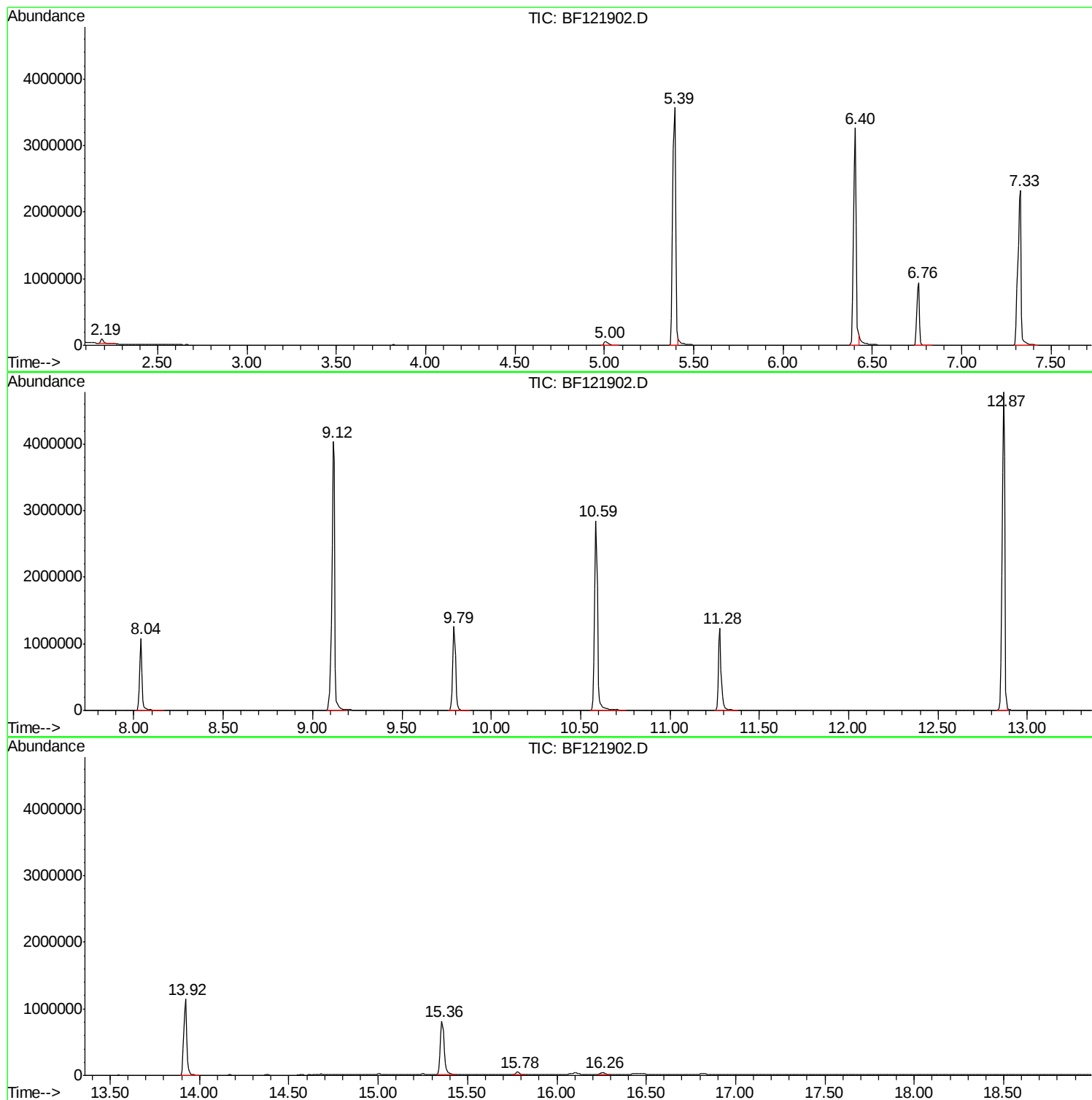
Sum of corrected areas: 30828607

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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF091020.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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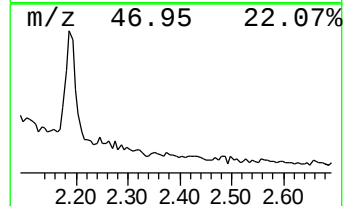
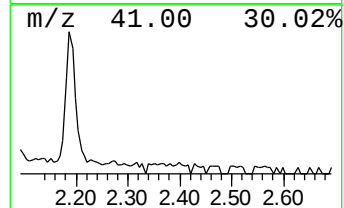
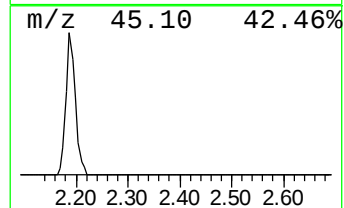
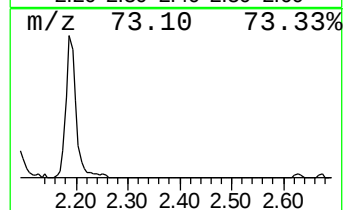
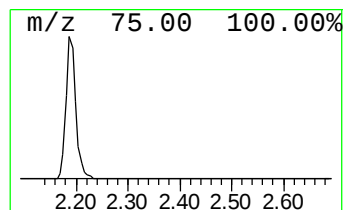
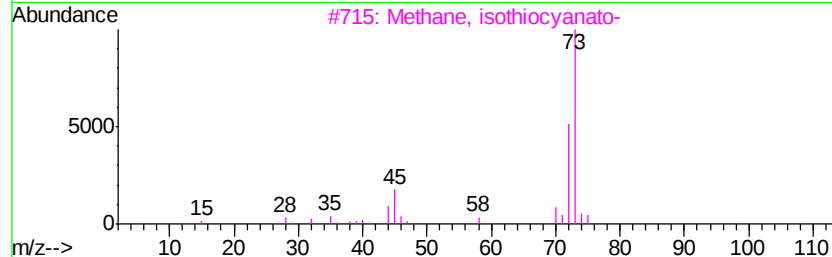
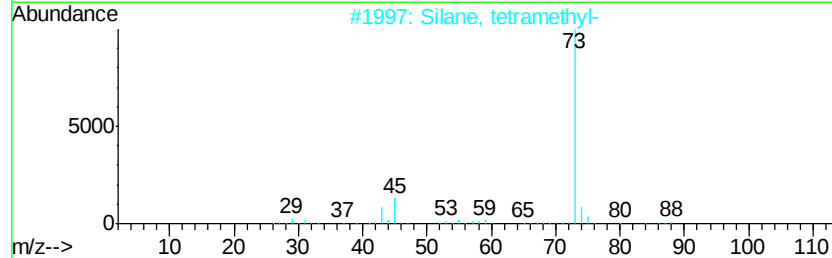
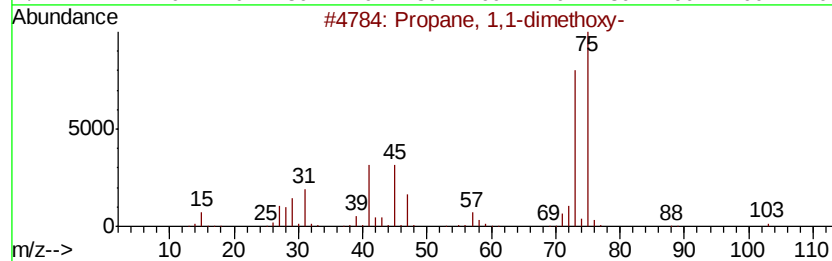
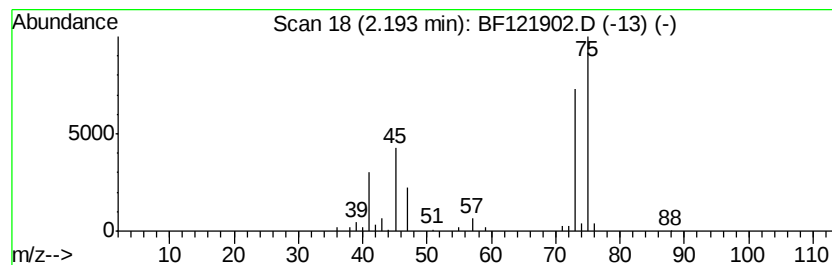
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TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.19	2.45 ng	103842	1,4-Dichlorobenzene-d4	6.76

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Propane, 1,1-dimethoxy-	104	C5H12O2	004744-10-9	72
2			Silane, tetramethyl-	88	C4H12Si	000075-76-3	9
3			Methane, isothiocyanato-	73	C2H3NS	000556-61-6	9
4			1,3-Dioxolane, 2-(1,1-dimethylet...	130	C7H14O2	002568-29-8	9
5			Pentane, 3-methoxy-	102	C6H14O	036839-67-5	8



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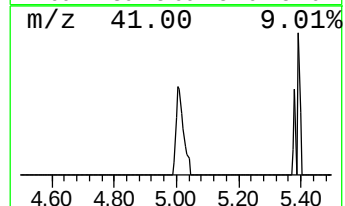
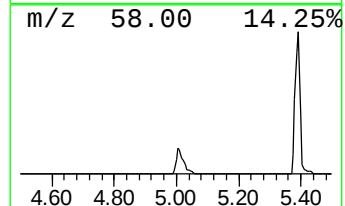
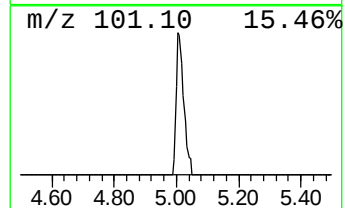
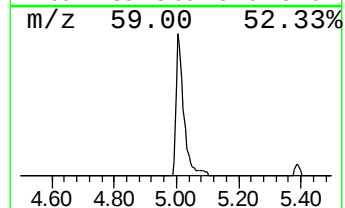
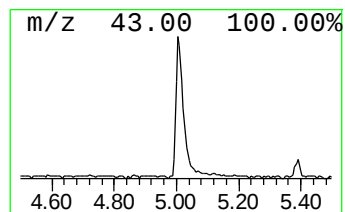
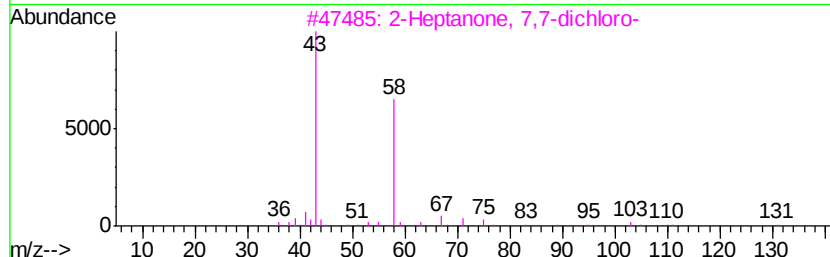
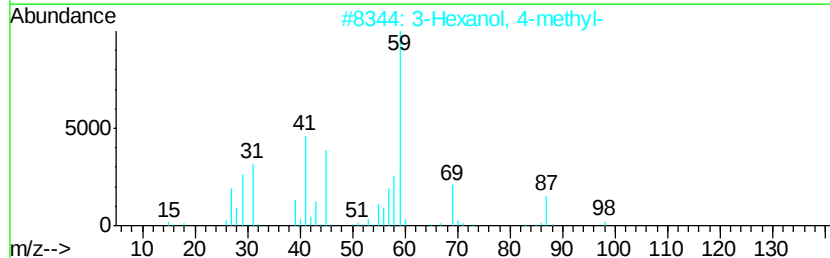
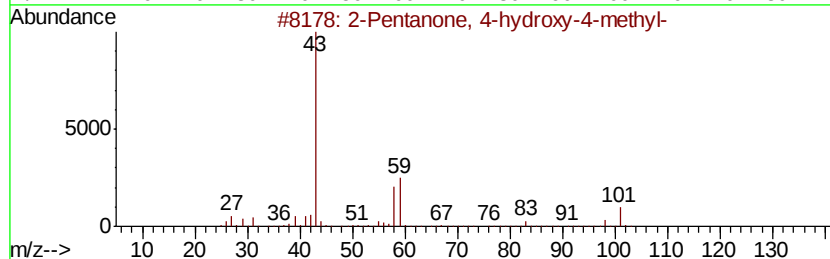
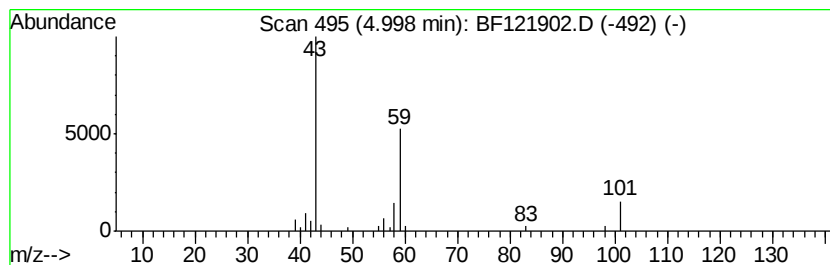
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TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.00	2.17 ng	92321	1,4-Dichlorobenzene-d4	6.76

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	53
2			3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	28
3			2-Heptanone, 7,7-dichloro-	182	C7H12Cl2O	066241-43-8	9
4			Butyl aldoxime, 3-methyl-, syn-	101	C5H11NO	005780-40-5	9
5			3-Octanol	130	C8H18O	000589-98-0	9



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TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Propane, 1,1-dime...	2.19	2.5	ng	103842	1	6.76	849282	20.0
2-Pentanone, 4-hy...	5.00	2.2	ng	92321	1	6.76	849282	20.0