

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP120219\
 Data File : BP001125.D
 Acq On : 02 Dec 2019 14:45
 Operator : JU
 Sample : PB125152BL
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 PB125152BL

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_P\METHODS\8270-BP112719.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	5.645	600	605	623	rBV	549440	918495	15.64%	2.134%
2	6.240	698	706	710	rBV	3296196	4474715	76.19%	10.395%
3	7.775	959	967	971	rBV	3629829	4711318	80.21%	10.945%
4	7.963	993	999	1004	rBV	3522134	4571141	77.83%	10.619%
5	8.328	1055	1061	1068	rBV	739227	875745	14.91%	2.034%
6	8.569	1096	1102	1106	rBV	2989992	3533921	60.17%	8.210%
7	9.210	1203	1211	1215	rBV	2307049	2976854	50.68%	6.916%
8	10.357	1401	1406	1412	rBV	911755	1106103	18.83%	2.570%
9	12.145	1702	1710	1714	rBV	4367988	5405058	92.03%	12.557%
10	13.222	1887	1893	1903	rBV	1047455	1307973	22.27%	3.039%
11	14.521	2106	2114	2136	rBV	2588416	3425760	58.33%	7.958%
12	15.621	2294	2301	2311	rBV	1188966	1374141	23.40%	3.192%
13	17.910	2684	2690	2694	rBV	5100205	5873442	100.00%	13.645%
14	19.262	2915	2920	2932	rBV	1142660	1270340	21.63%	2.951%
15	21.045	3217	3223	3236	rVB	917533	1220560	20.78%	2.836%

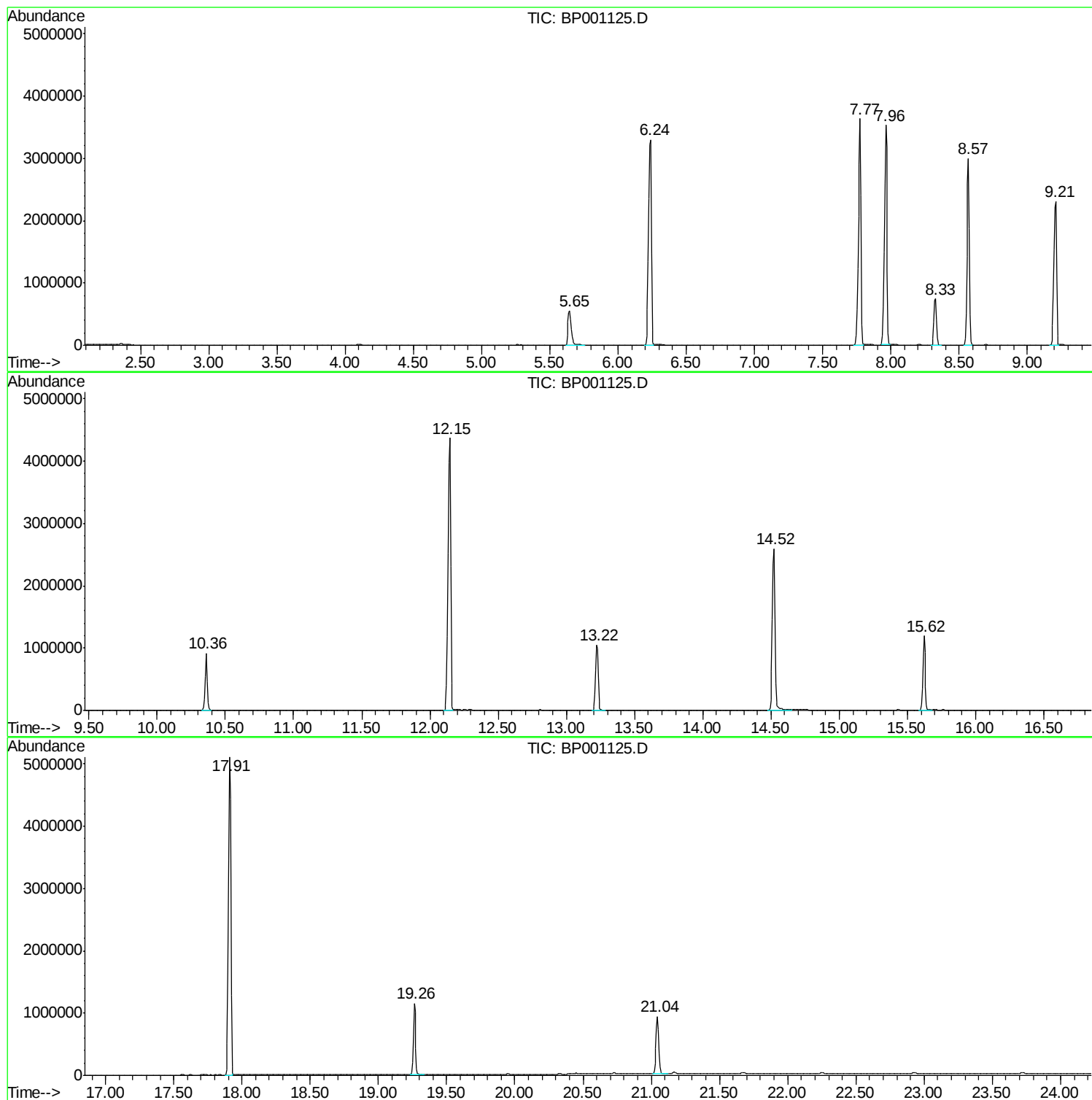
Sum of corrected areas: 43045566

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Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP112719.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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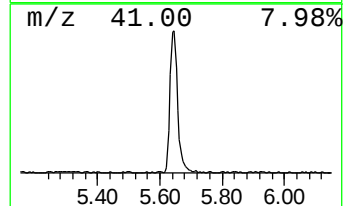
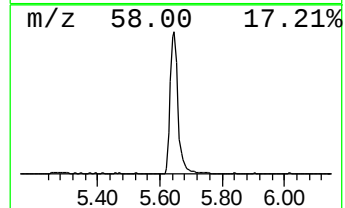
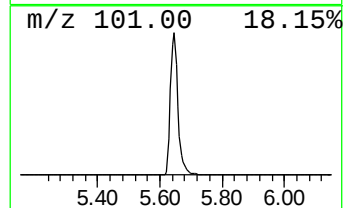
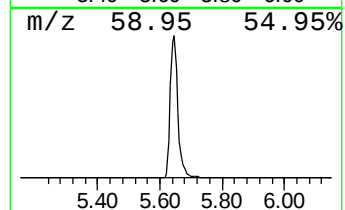
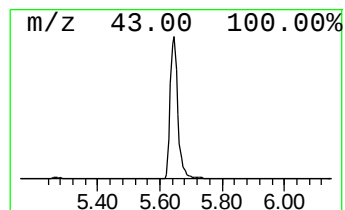
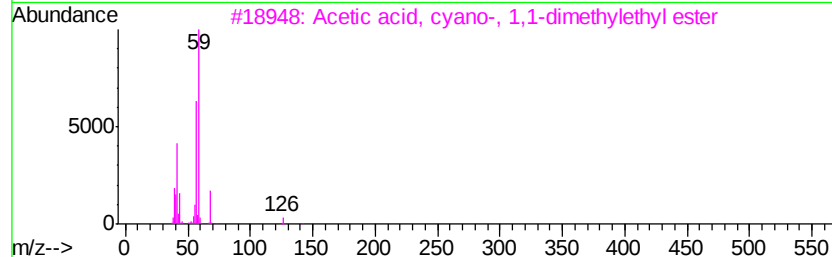
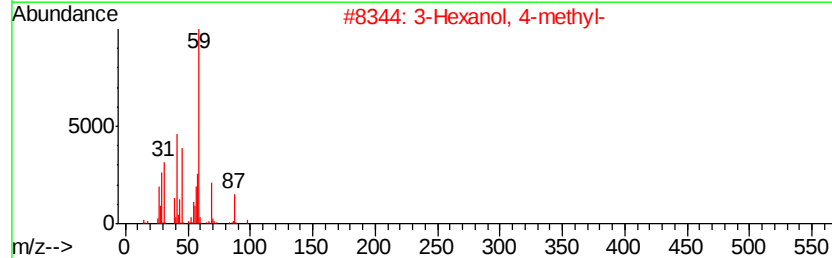
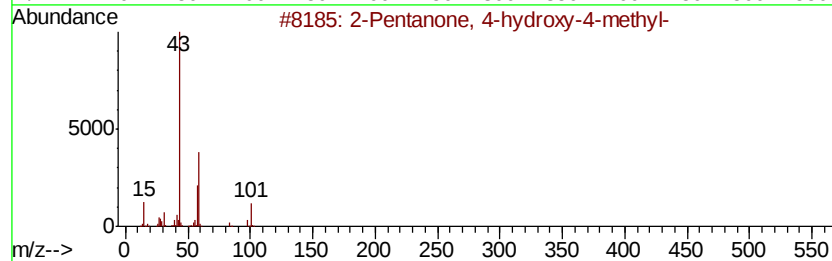
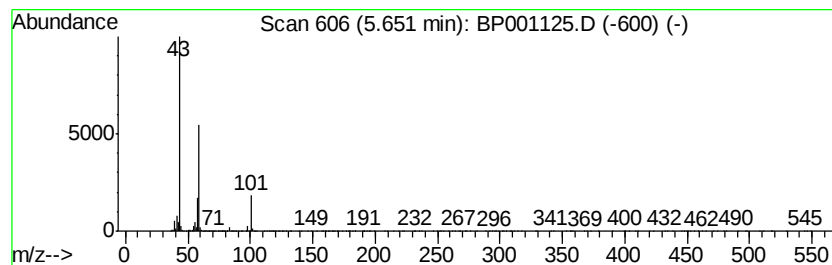
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.65	20.98 ng	918495	1,4-Dichlorobenzene-d4	8.33

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	59
2			3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	33
3			Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	25
4			1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	10
5			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	10



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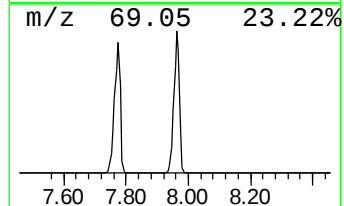
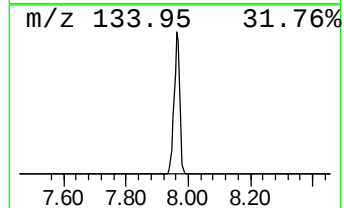
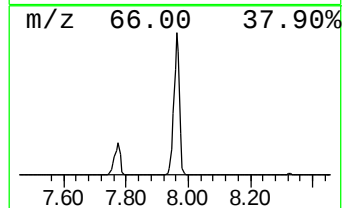
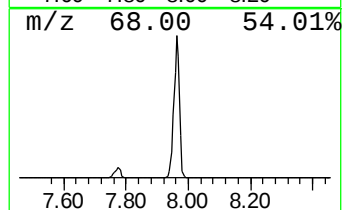
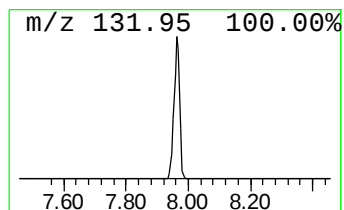
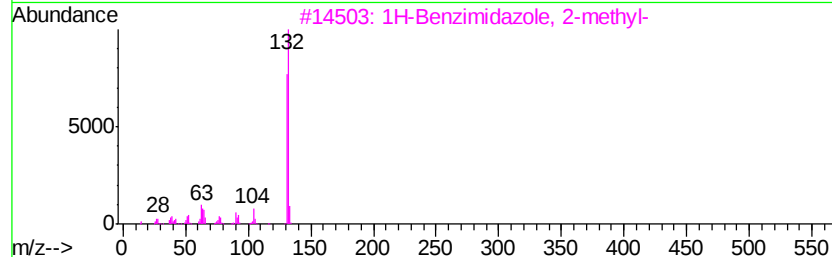
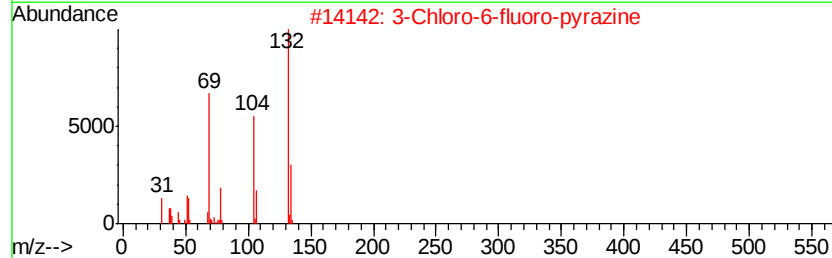
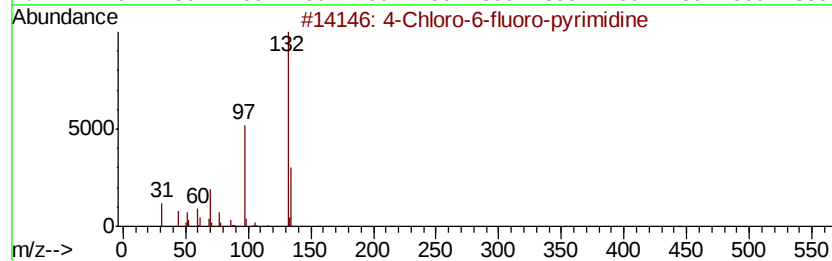
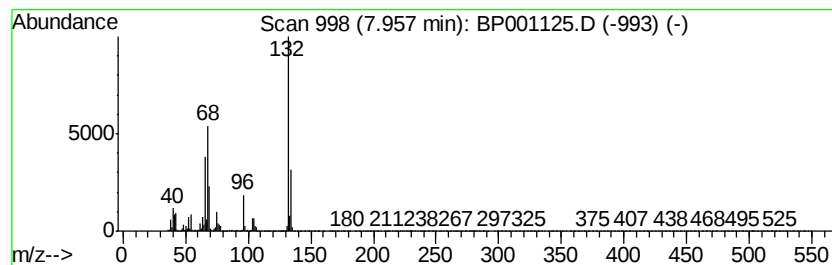
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Peak Number 2 unknown7.96 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.96	104.39 ng	4571140	1,4-Dichlorobenzene-d4	8.33

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	4-Chloro-6-fluoro-pyrimidine	132	C4H2ClFN2	051422-01-6	27
2		3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
3		1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	22
4		1,3-Cyclopentanedione, 2-chloro-	132	C5H5ClO2	014203-19-1	13
5		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	12



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.65	21.0	ng	918495	1	8.33	875745	20.0
unknown7.96	7.96	104.4	ng	4571140	1	8.33	875745	20.0