

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF012419\
 Data File : BF112207.D
 Acq On : 24 Jan 2019 14:23
 Operator : JU/SJ
 Sample : PB116593BL
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB116593BL

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-BF011619.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	5.098	507	512	523	rBV	215933	308585	3.46%	0.490%
2	5.498	576	580	597	rBV	5582958	5788646	64.89%	9.190%
3	6.498	745	750	767	rBV	5036236	6088396	68.25%	9.665%
4	6.628	767	772	776	rBV	5920056	6350526	71.19%	10.082%
5	6.851	806	810	818	rBV	1689754	1434896	16.09%	2.278%
6	7.010	832	837	840	rBV	5477705	5062278	56.75%	8.036%
7	7.416	900	906	909	rBV	3317020	3949298	44.27%	6.270%
8	8.133	1022	1028	1044	rBV	1894584	1918764	21.51%	3.046%
9	9.210	1204	1211	1214	rBV	7530008	8125140	91.09%	12.899%
10	9.886	1320	1326	1344	rBV	2214604	2331435	26.14%	3.701%
11	10.680	1455	1461	1483	rBV	4874674	5551961	62.24%	8.814%
12	11.374	1574	1579	1587	rBV	2756104	2553422	28.63%	4.054%
13	12.963	1842	1849	1852	rBV	8078490	8920217	100.00%	14.161%
14	14.021	2024	2029	2037	rBV	2376859	2443703	27.40%	3.879%
15	15.492	2273	2279	2292	rBV	1498111	2164513	24.27%	3.436%

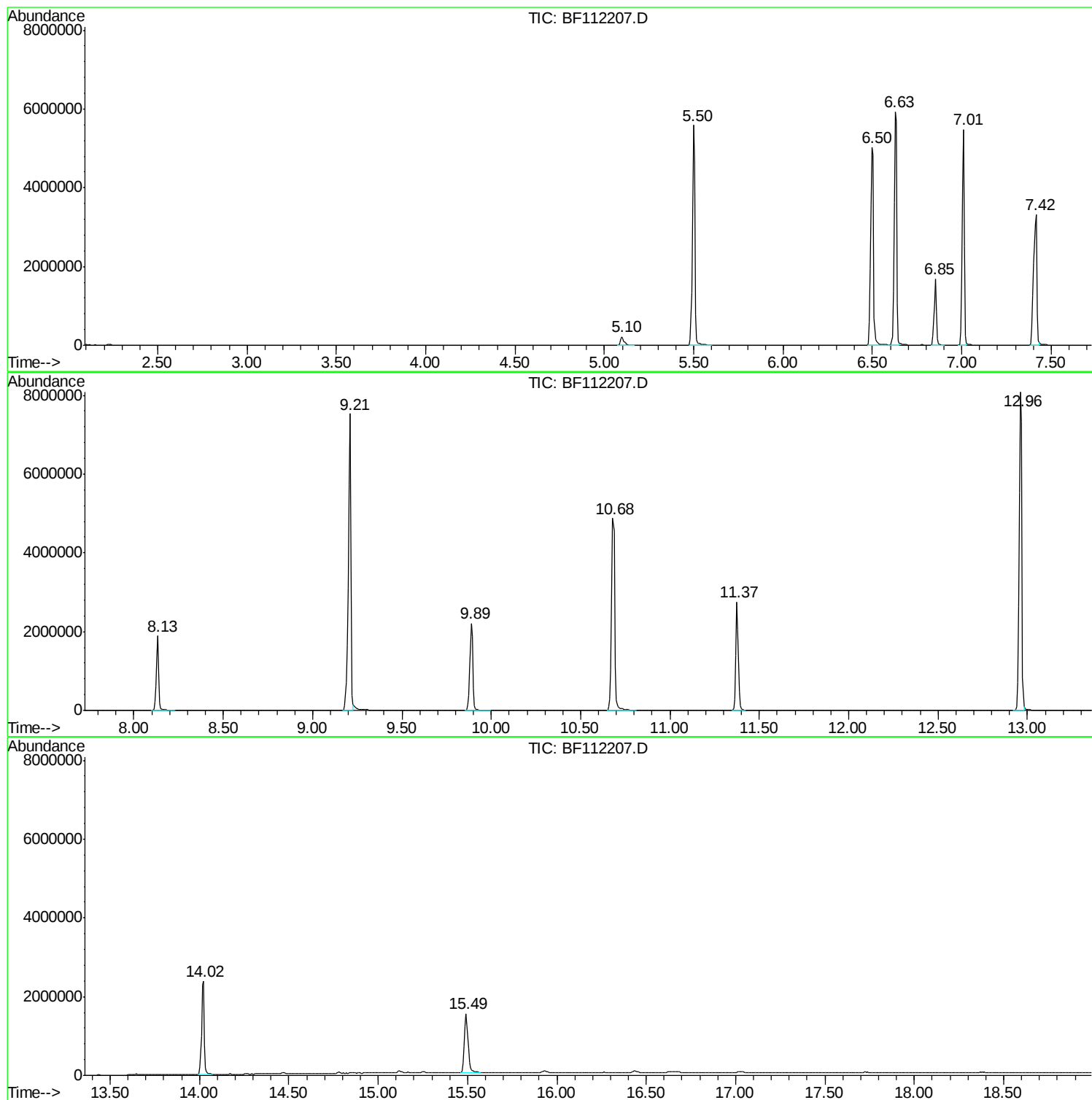
Sum of corrected areas: 62991780

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



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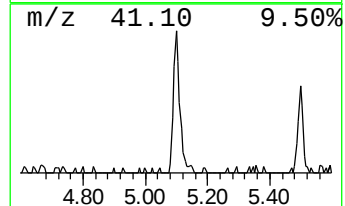
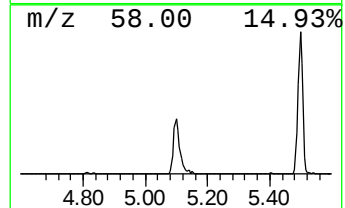
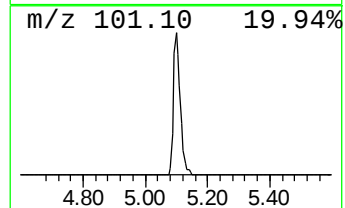
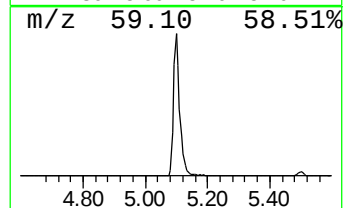
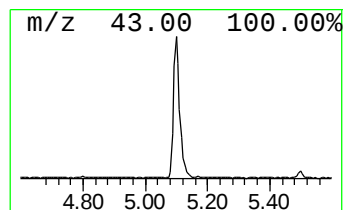
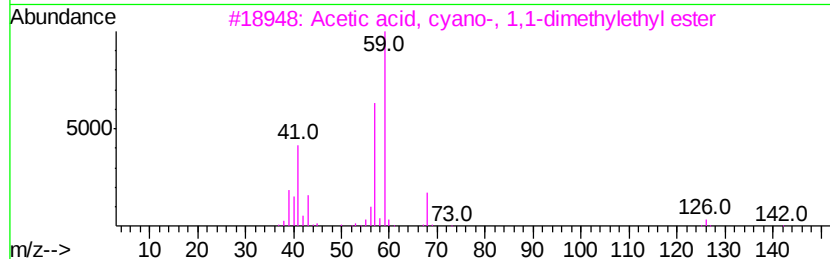
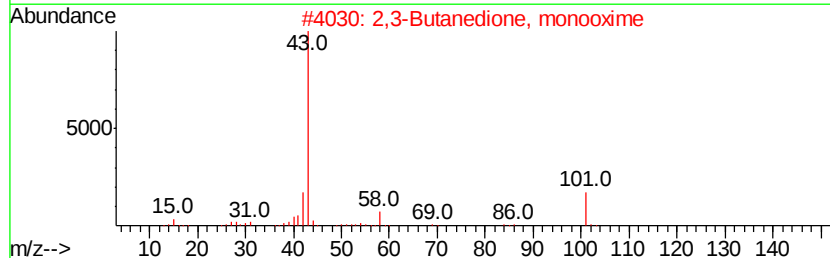
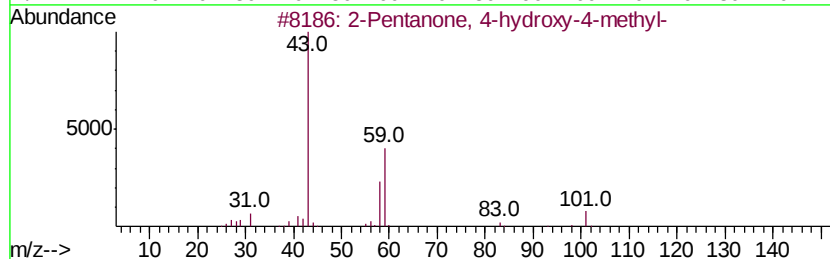
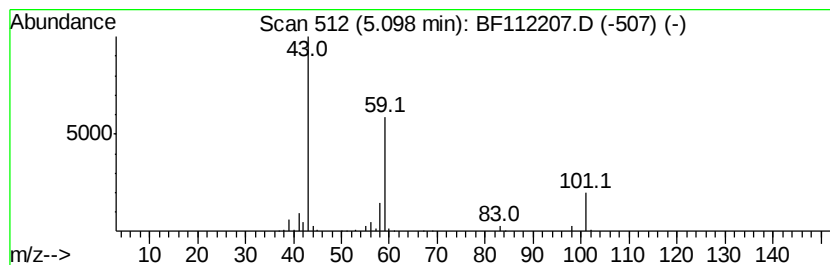
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.	
5.10	4.30 ng	308585	1,4-Dichlorobenzene-d4	6.85	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	42
2	2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	35
3	Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	23
4	Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9
5	3-Hexanol	102	C6H14O	000623-37-0	9



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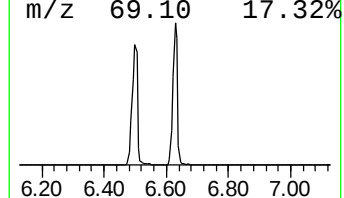
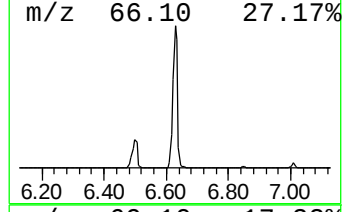
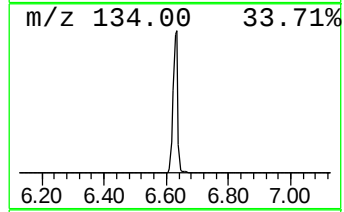
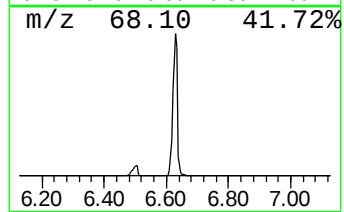
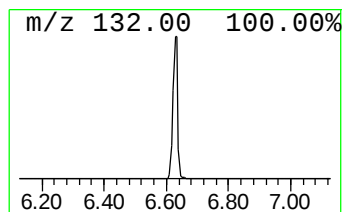
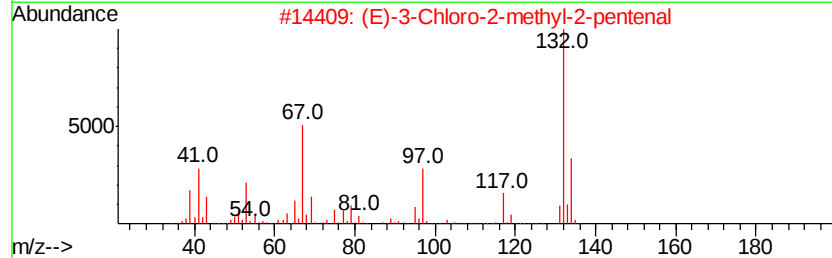
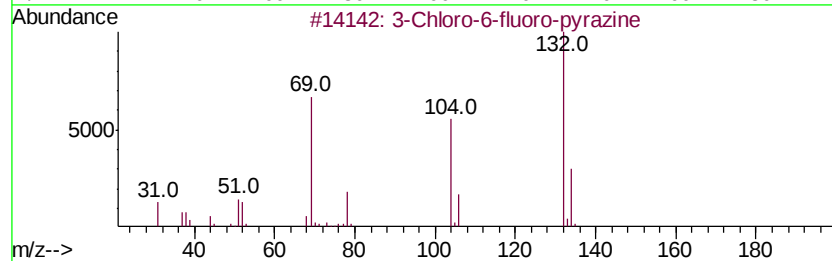
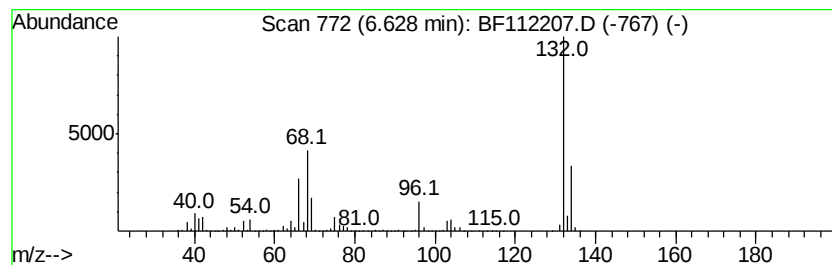
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Peak Number 2 unknown6.63 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.63	88.52 ng	6350530	1,4-Dichlorobenzene-d4	6.85

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	17
3		Tranvlcypropine-propionyl	189	C12H15NO	1000123-86-3	10
4		Benzene, 1-ethenyl-3,5-dimethyl-	132	C10H12	005379-20-4	9
5		4-Chloro-2-fluoro-pyrimidine	132	C4H2ClFN2	051422-00-5	9



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TIC Top Hit name	RT	EstConc	Units	Response	---Internal Standard---			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.10	4.3	ng	308585	1	6.85	1434900	20.0
unknown6.63	6.63	88.5	ng	6350530	1	6.85	1434900	20.0