

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF073018\
 Data File : BF107804.D
 Acq On : 30 Jul 2018 17:13
 Operator : JU/SJ
 Sample : PB111493BL
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB111493BL

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-BF073018.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	5.201	483	487	499	rBV	412691	570837	11.81%	1.638%
2	5.601	551	555	578	rBV	2094532	2982469	61.71%	8.557%
3	6.601	721	725	744	rBV	1845893	3077270	63.67%	8.829%
4	6.731	744	747	769	rVB	2491383	3403676	70.43%	9.766%
5	6.960	783	786	808	rVB	529621	874588	18.10%	2.509%
6	7.113	808	812	838	rVB	2435780	2706445	56.00%	7.765%
7	7.525	878	882	917	rBV	1262099	2220011	45.94%	6.370%
8	8.242	1000	1004	1030	rBV	828156	1222588	25.30%	3.508%
9	9.313	1181	1186	1222	rBV	3477984	4312361	89.23%	12.373%
10	10.001	1298	1303	1322	rBV	1133042	1409017	29.15%	4.043%
11	10.795	1434	1438	1467	rBV	2059533	2894866	59.90%	8.306%
12	11.495	1553	1557	1585	rBV	1192268	1629498	33.72%	4.675%
13	13.077	1822	1826	1852	rBV	3994168	4832878	100.00%	13.866%
14	14.148	2004	2008	2025	rBV	1098642	1467078	30.36%	4.209%
15	15.677	2259	2268	2286	rBV	581342	1190819	24.64%	3.417%
16	16.101	2336	2340	2347	rBV3	31596	58992	1.22%	0.169%

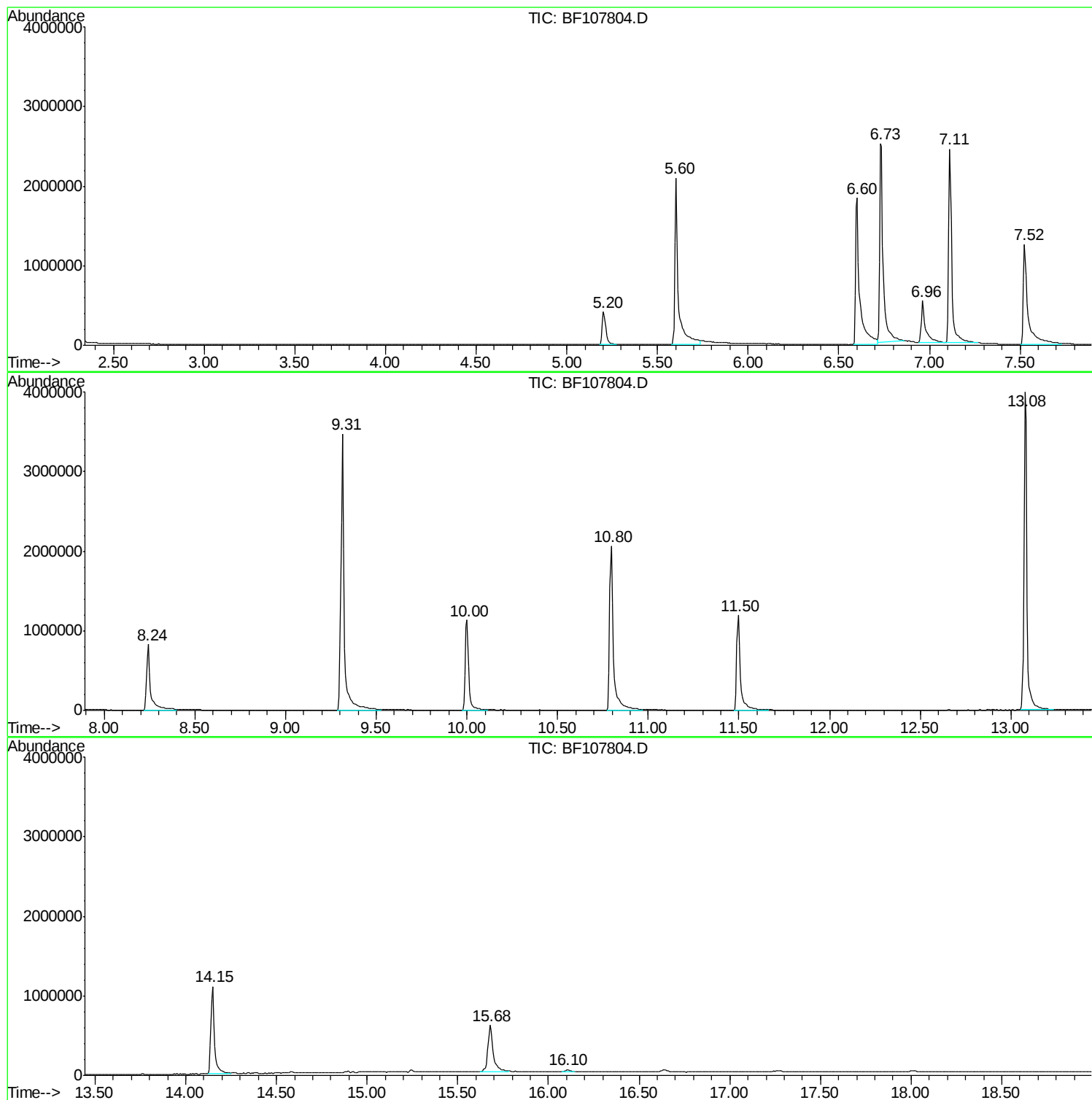
Sum of corrected areas: 34853393

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



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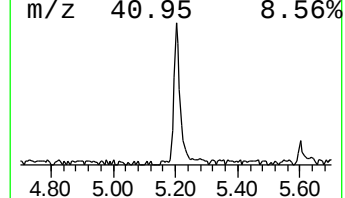
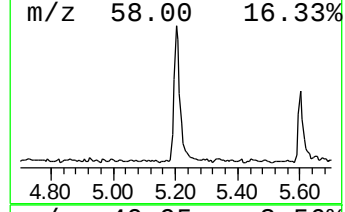
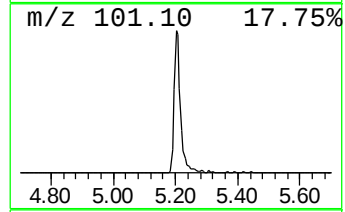
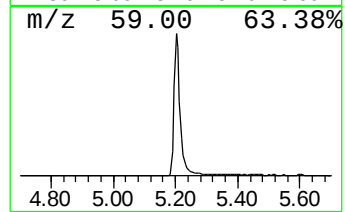
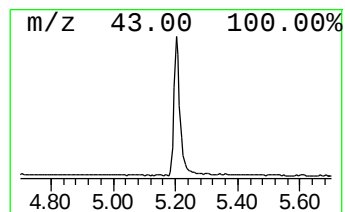
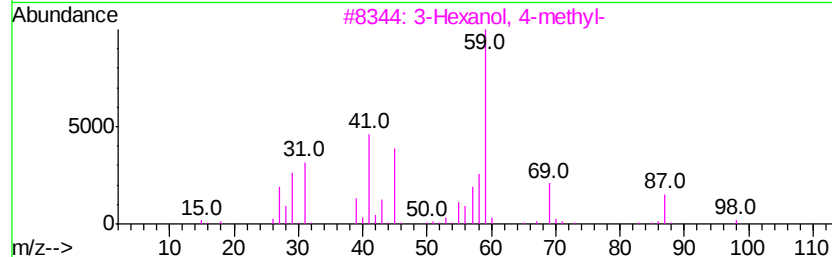
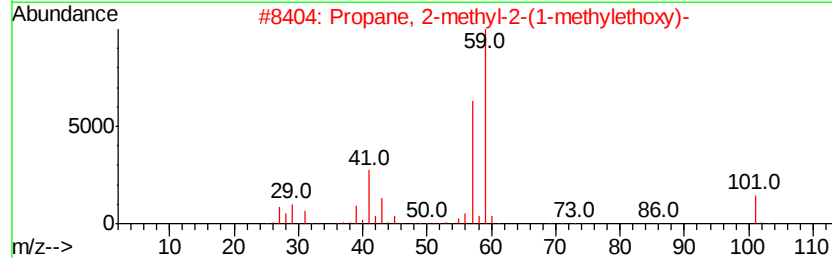
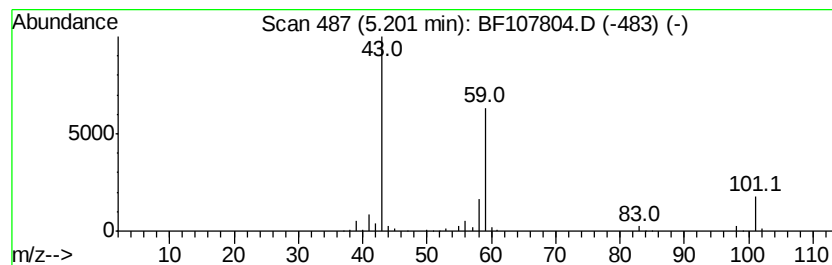
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.20	13.05 ng	570837	1,4-Dichlorobenzene-d4	6.96

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	59
2		Propane, 2-methyl-2-(1-methyleth...	116	C7H16O	017348-59-3	36
3		3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	33
4		2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	33
5		Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	28



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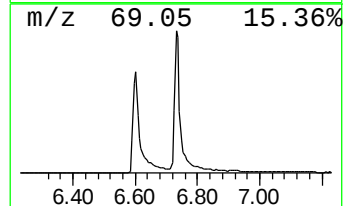
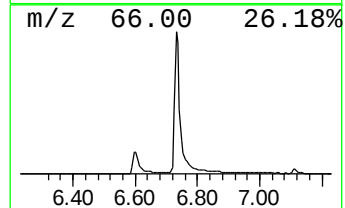
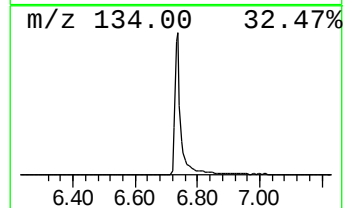
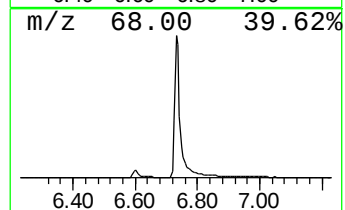
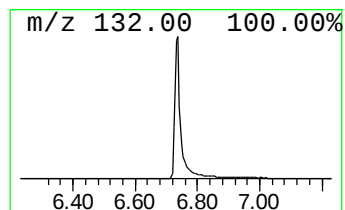
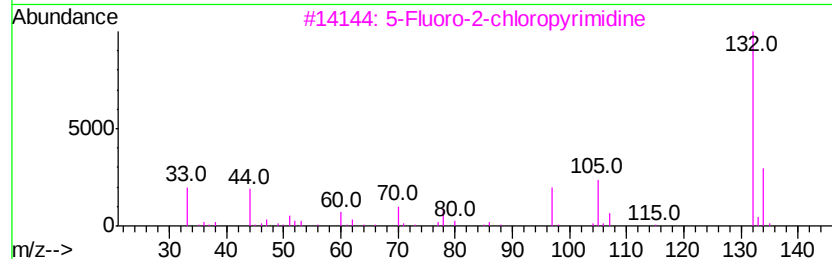
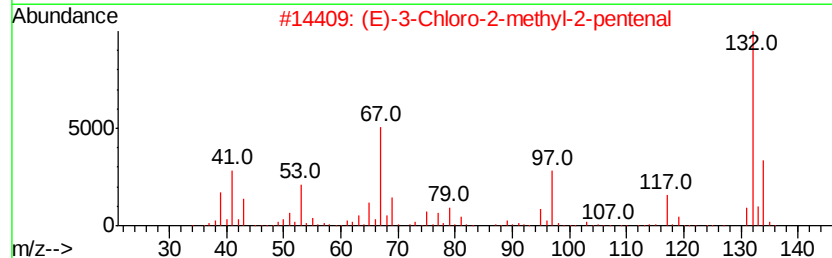
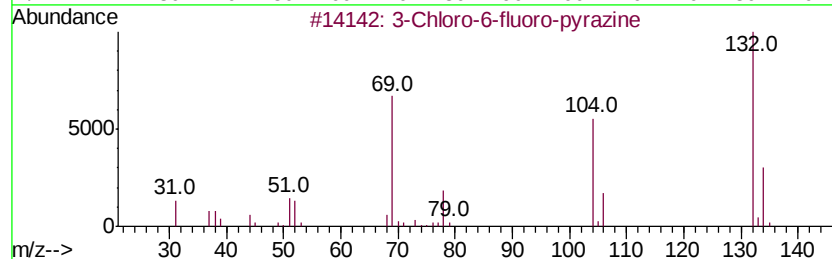
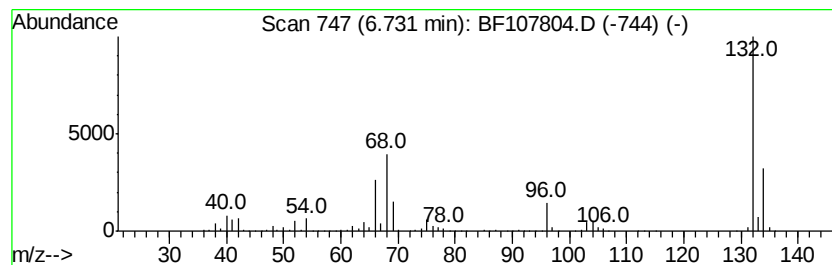
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Peak Number 2 unknown6.73 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.73	77.84 ng	3403680	1,4-Dichlorobenzene-d4	6.96

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		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	16
4		5-Aminoindole	132	C8H8N2	005192-03-0	12
5		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.20	13.1	ng	570837	1	6.96	874588	20.0
unknown6.73	6.73	77.8	ng	3403680	1	6.96	874588	20.0