

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF110518\
Data File : BF110486.D
Acq On : 6 Nov 2018 1:46
Operator : JU/SJ
Sample : PB114474BL
Misc :
ALS Vial : 31 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
PB114474BL

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-BF102318.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	2.346	39	44	49	rVB	24073	34197	1.42%	0.227%
2	5.198	525	529	542	rBV	60554	83578	3.48%	0.555%
3	5.581	590	594	609	rBV	1441413	1292992	53.82%	8.580%
4	6.575	759	763	777	rBV	1410692	1352309	56.29%	8.973%
5	6.722	784	788	797	rBV	1708048	1417735	59.01%	9.407%
6	6.951	824	827	837	rBV	650147	545095	22.69%	3.617%
7	7.110	850	854	869	rBV	1826339	1648385	68.61%	10.938%
8	7.516	918	923	944	rBV	1199481	1307549	54.42%	8.676%
9	8.239	1041	1046	1066	rBV	571870	713745	29.71%	4.736%
10	9.310	1222	1228	1244	rBV	2679192	2402594	100.00%	15.942%
11	9.998	1341	1345	1359	rVB	638243	725340	30.19%	4.813%
12	10.786	1475	1479	1494	rBV	587160	645129	26.85%	4.281%
13	11.486	1594	1598	1610	rBV	505239	588503	24.49%	3.905%
14	13.068	1863	1867	1884	rVB	1741044	1564386	65.11%	10.380%
15	14.133	2044	2048	2060	rBV	374184	437552	18.21%	2.903%
16	15.662	2303	2308	2317	rVB	200100	311526	12.97%	2.067%

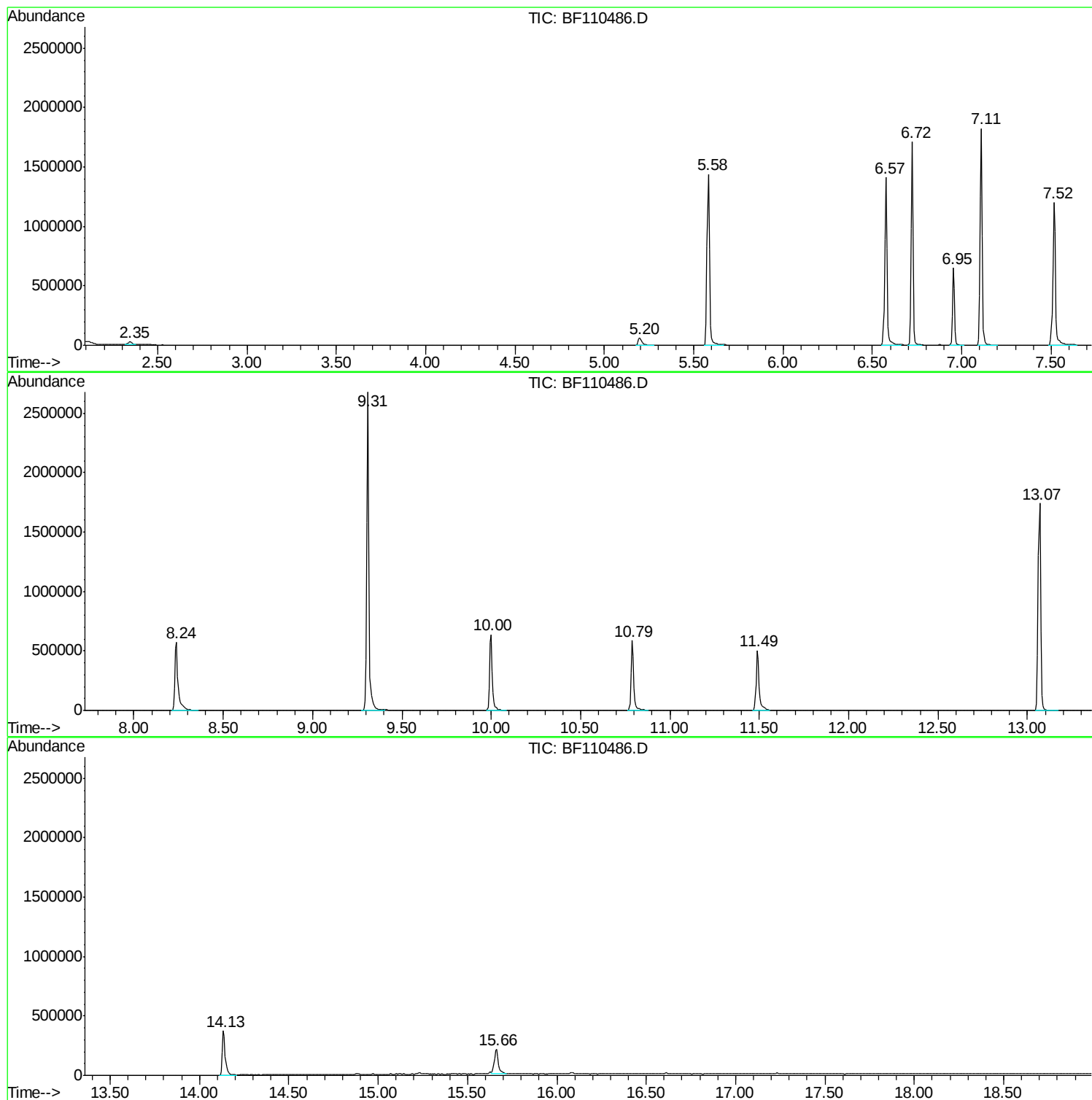
Sum of corrected areas: 15070615

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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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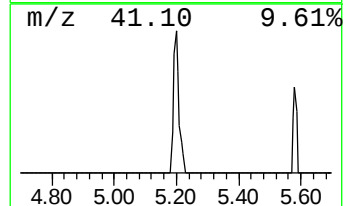
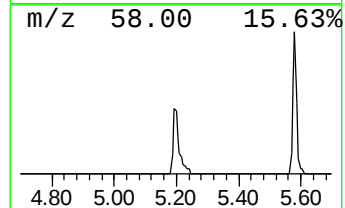
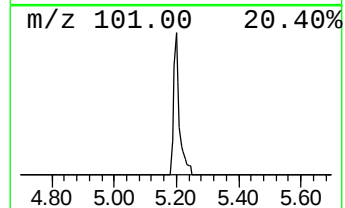
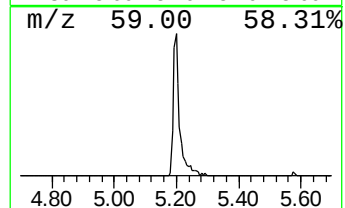
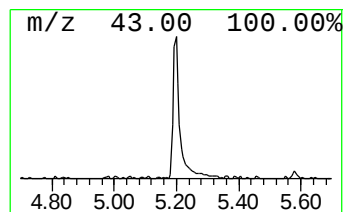
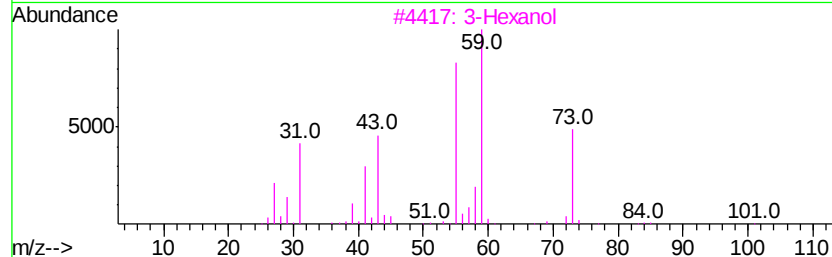
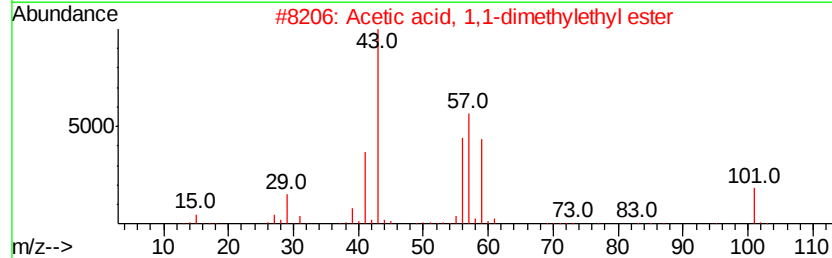
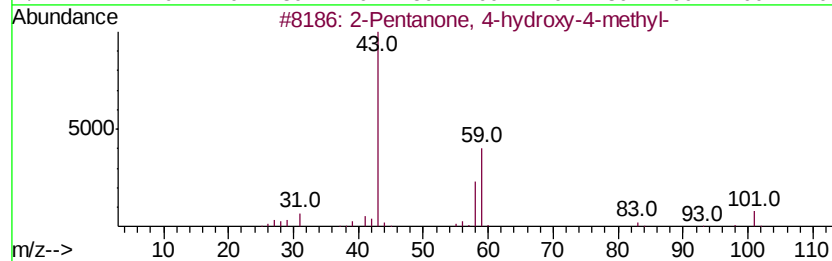
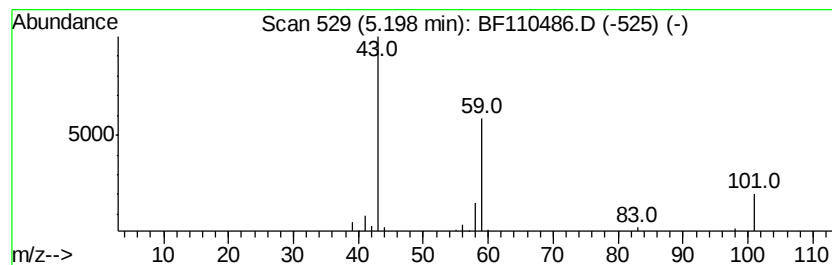
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TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.20	3.07 ng	83578	1,4-Dichlorobenzene-d4	6.95

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	42
2			Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	39
3			3-Hexanol	102	C6H14O	000623-37-0	28
4			5-Hexen-2-one	98	C6H10O	000109-49-9	9
5			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9



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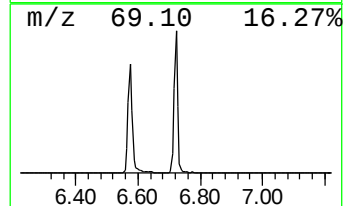
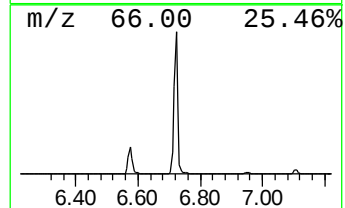
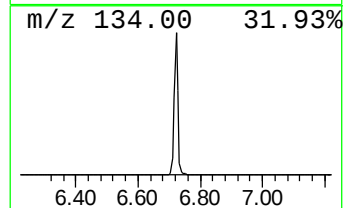
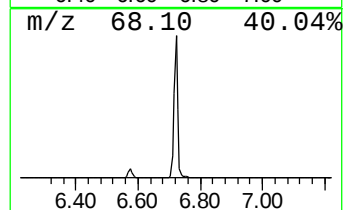
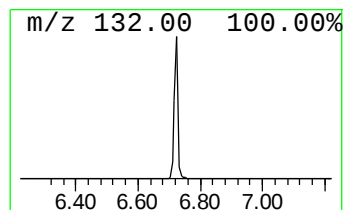
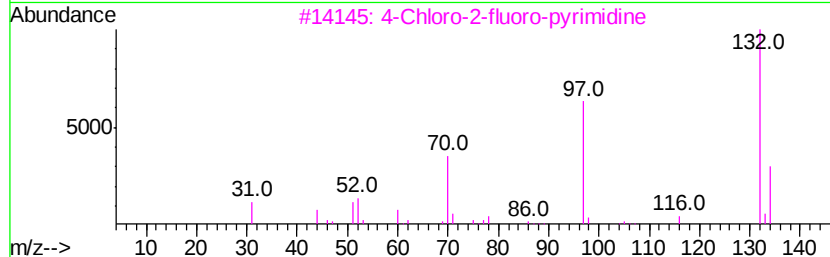
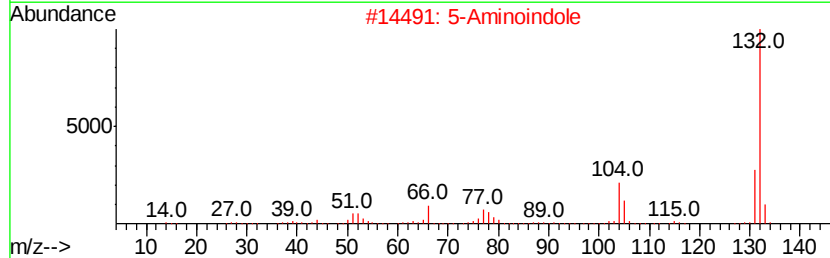
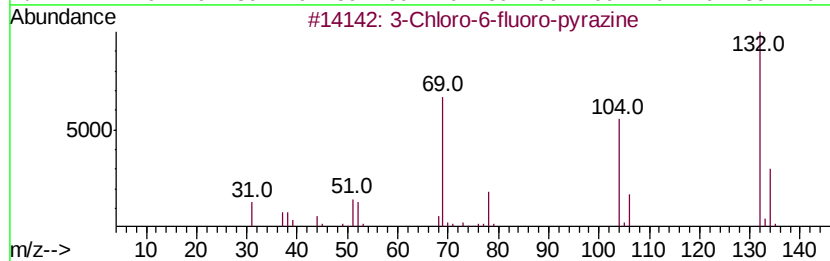
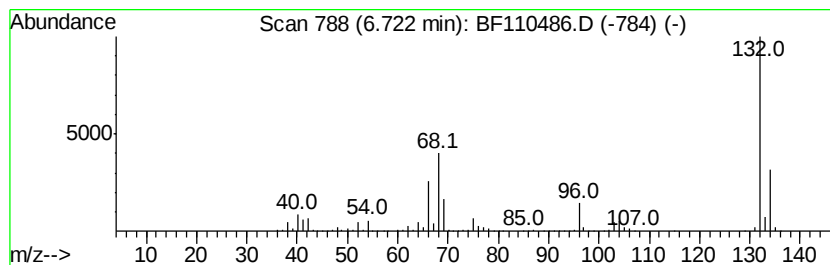
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Peak Number 2 unknown6.72 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.72	52.02 ng	1417740	1,4-Dichlorobenzene-d4	6.95

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2			5-Aminoindole	132	C8H8N2	005192-03-0	12
3			4-Chloro-2-fluoro-pyrimidine	132	C4H2ClFN2	051422-00-5	9
4			(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	9
5			5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.20	3.1	ng	83578	1	6.95	545095	20.0
unknown6.72	6.72	52.0	ng	1417740	1	6.95	545095	20.0