

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA_F\DATA\BF110717\
 Data File : BF100279.D
 Acq On : 7 Nov 2017 12:56
 Operator : SJ/JU
 Sample : I6336-01
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 CF-1B(1-2)

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:\HPCHEM1\BNA_F\METHODS\8270-BF102317.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	4.934	482	484	500	rBV	59399	129125	4.26%	0.692%
2	5.340	550	553	583	rBV	549834	1048493	34.60%	5.616%
3	6.369	726	728	744	rBV	387078	659155	21.75%	3.530%
4	6.487	744	748	763	rVB	1577177	1897336	62.61%	10.162%
5	6.710	783	786	798	rBV	520972	534164	17.63%	2.861%
6	6.863	808	812	835	rBV	2034403	2002032	66.07%	10.723%
7	7.287	880	884	904	rBV	1368370	1504980	49.66%	8.061%
8	7.992	1001	1004	1017	rBV	754317	779216	25.71%	4.173%
9	9.063	1182	1186	1200	rVB	2994780	3030375	100.00%	16.230%
10	9.739	1298	1301	1314	rVB	852302	884009	29.17%	4.735%
11	10.539	1433	1437	1458	rBV	1436911	1464926	48.34%	7.846%
12	11.233	1552	1555	1567	rBV	956038	911016	30.06%	4.879%
13	12.816	1820	1824	1828	rBV	2559942	2754689	90.90%	14.754%
14	13.886	2003	2006	2021	rBV	600254	652516	21.53%	3.495%
15	15.333	2248	2252	2267	rBV	246251	418900	13.82%	2.244%

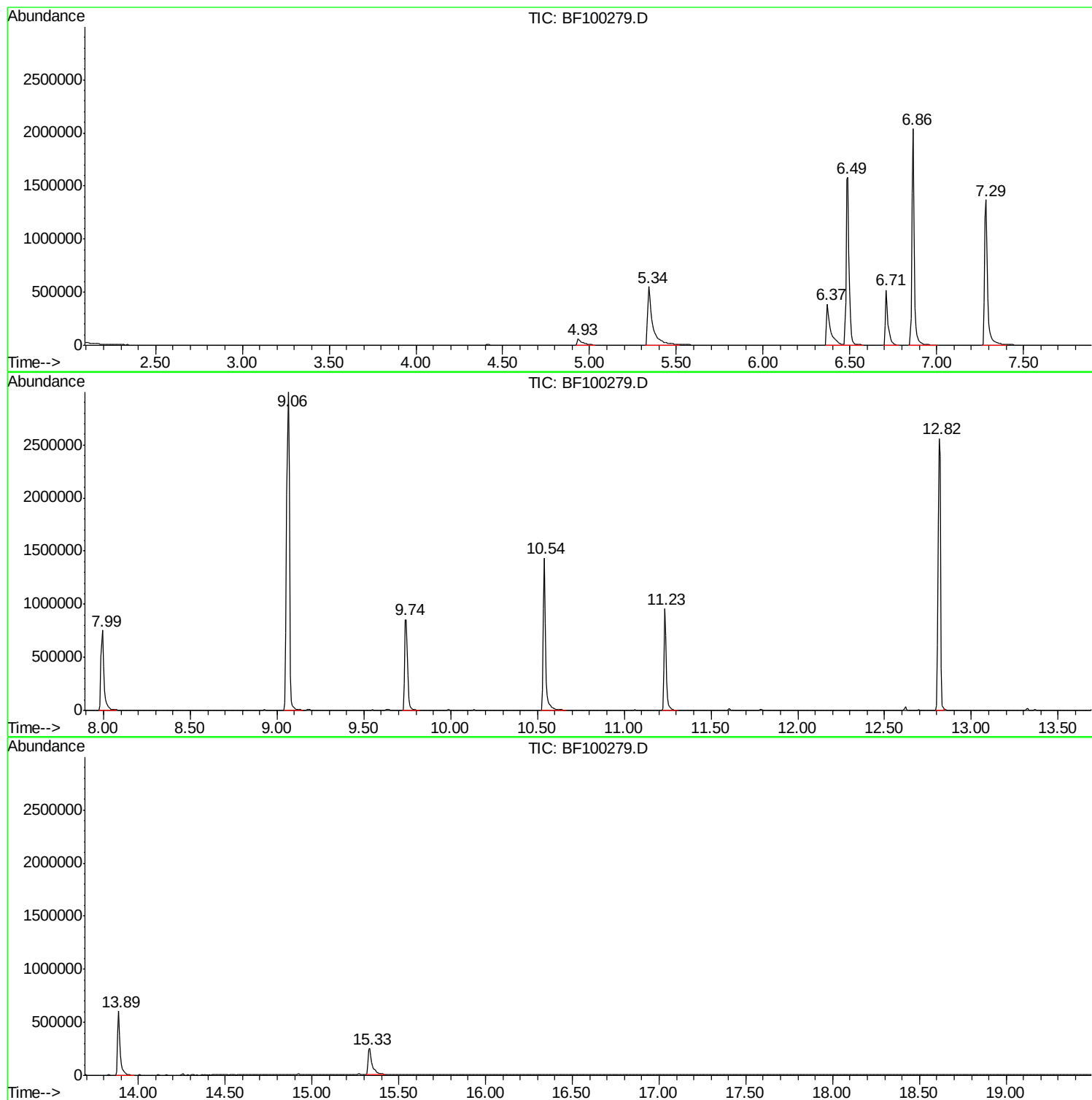
Sum of corrected areas: 18670932

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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF102317.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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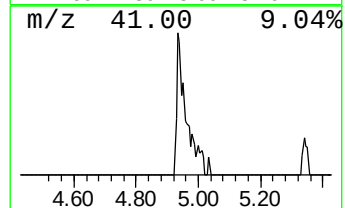
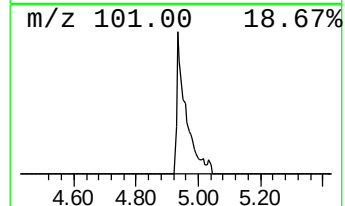
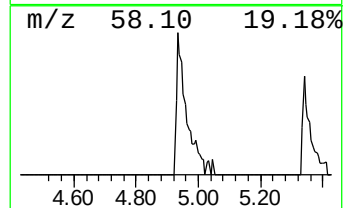
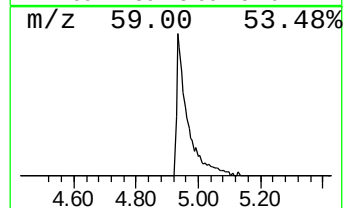
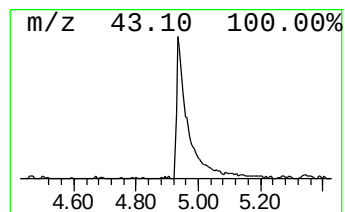
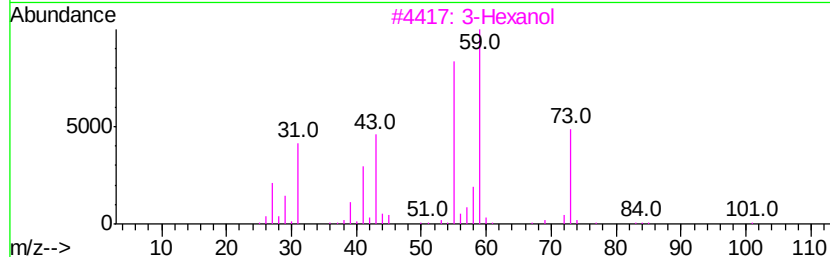
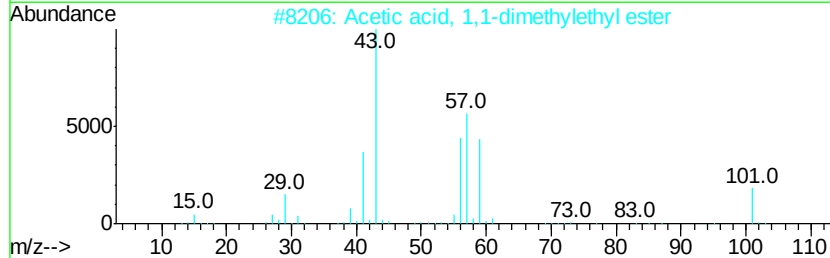
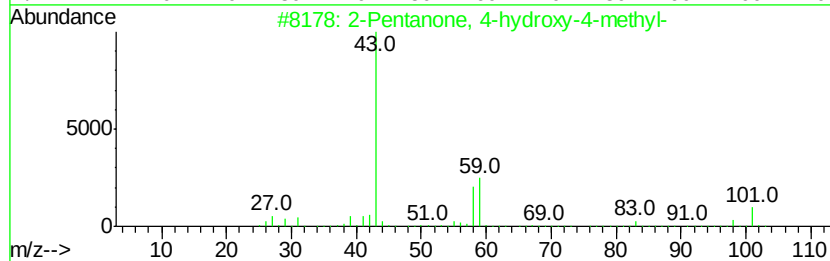
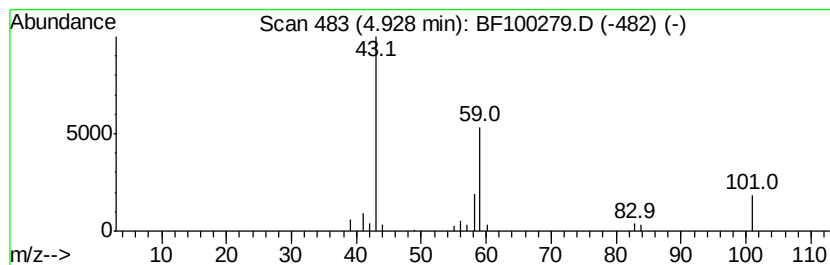
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.93	4.83 ng	129125	1,4-Dichlorobenzene-d4	6.71

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	45
2			Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	39
3			3-Hexanol	102	C6H14O	000623-37-0	28
4			3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	28
5			Butyl aldoxime, 3-methyl-, syn-	101	C5H11NO	005780-40-5	25



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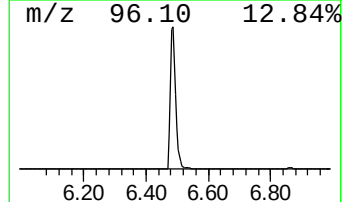
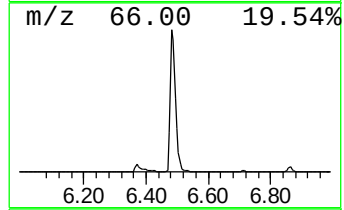
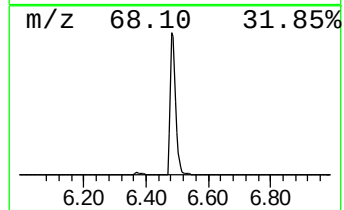
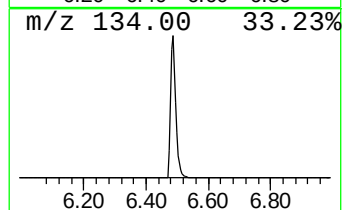
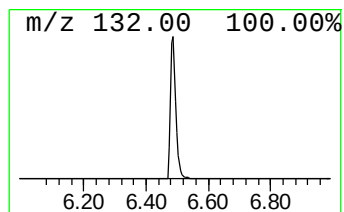
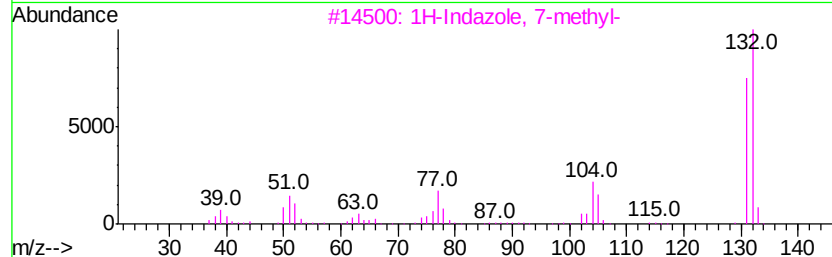
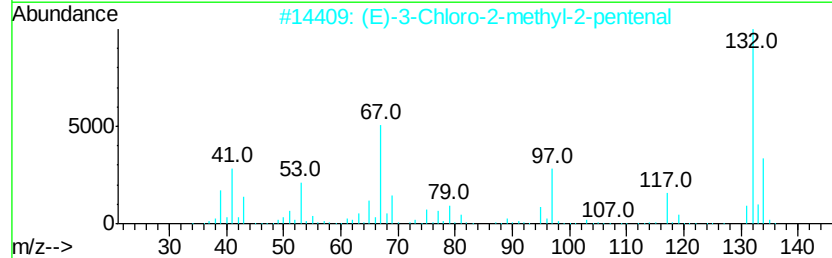
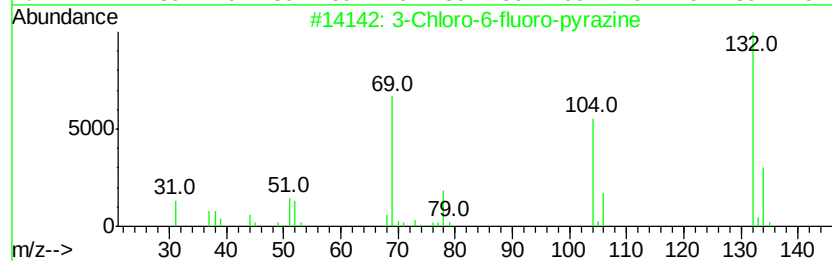
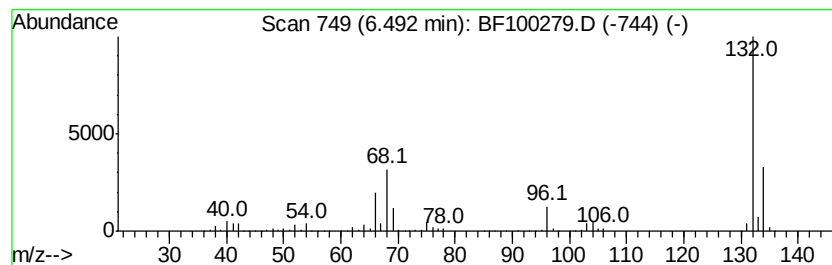
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Peak Number 2 3-Chloro-6-fluoro-pyrazine Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.49	71.04 ng	1897340	1,4-Dichlorobenzene-d4	6.71

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	52
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	33
3		1H-Indazole, 7-methyl-	132	C8H8N2	1000316-00-7	25
4		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	9
5		Xenon	132	Xe	007440-63-3	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	4.93	4.8	ng	129125	1	6.71	534164	20.0
3-Chloro-6-fluoro...	6.49	71.0	ng	1897340	1	6.71	534164	20.0