

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF091417\
 Data File : BF098534.D
 Acq On : 14 Sep 2017 18:06
 Operator : SJ/JU
 Sample : PB102300BL
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB102300BL

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-BF091117.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	4.822	462	465	476	rBV	156764	227444	6.48%	0.837%
2	5.251	534	538	561	rBV	2569659	2814814	80.23%	10.361%
3	6.298	711	716	732	rBV	2551981	2688993	76.64%	9.898%
4	6.416	732	736	740	rBV	3165377	2756918	78.58%	10.148%
5	6.645	771	775	781	rBV	835644	710272	20.24%	2.614%
6	6.798	797	801	812	rBV	2333813	2123562	60.53%	7.816%
7	7.216	867	872	884	rBV	1664949	1709065	48.71%	6.291%
8	7.928	989	993	1002	rBV	1102907	968068	27.59%	3.563%
9	9.004	1171	1176	1180	rBV	3370131	3137462	89.42%	11.548%
10	9.680	1286	1291	1297	rBV	1135258	1092686	31.14%	4.022%
11	10.469	1421	1425	1444	rVB	2050669	1861539	53.06%	6.852%
12	11.163	1539	1543	1546	rBV	1273595	1192681	33.99%	4.390%
13	12.745	1807	1812	1815	rBV	3709621	3508500	100.00%	12.914%
14	13.792	1986	1990	1994	rBV	1344971	1189073	33.89%	4.377%
15	15.192	2222	2228	2240	rBV	908663	1149003	32.75%	4.229%
16	16.045	2369	2373	2376	rBV6	24731	37886	1.08%	0.139%

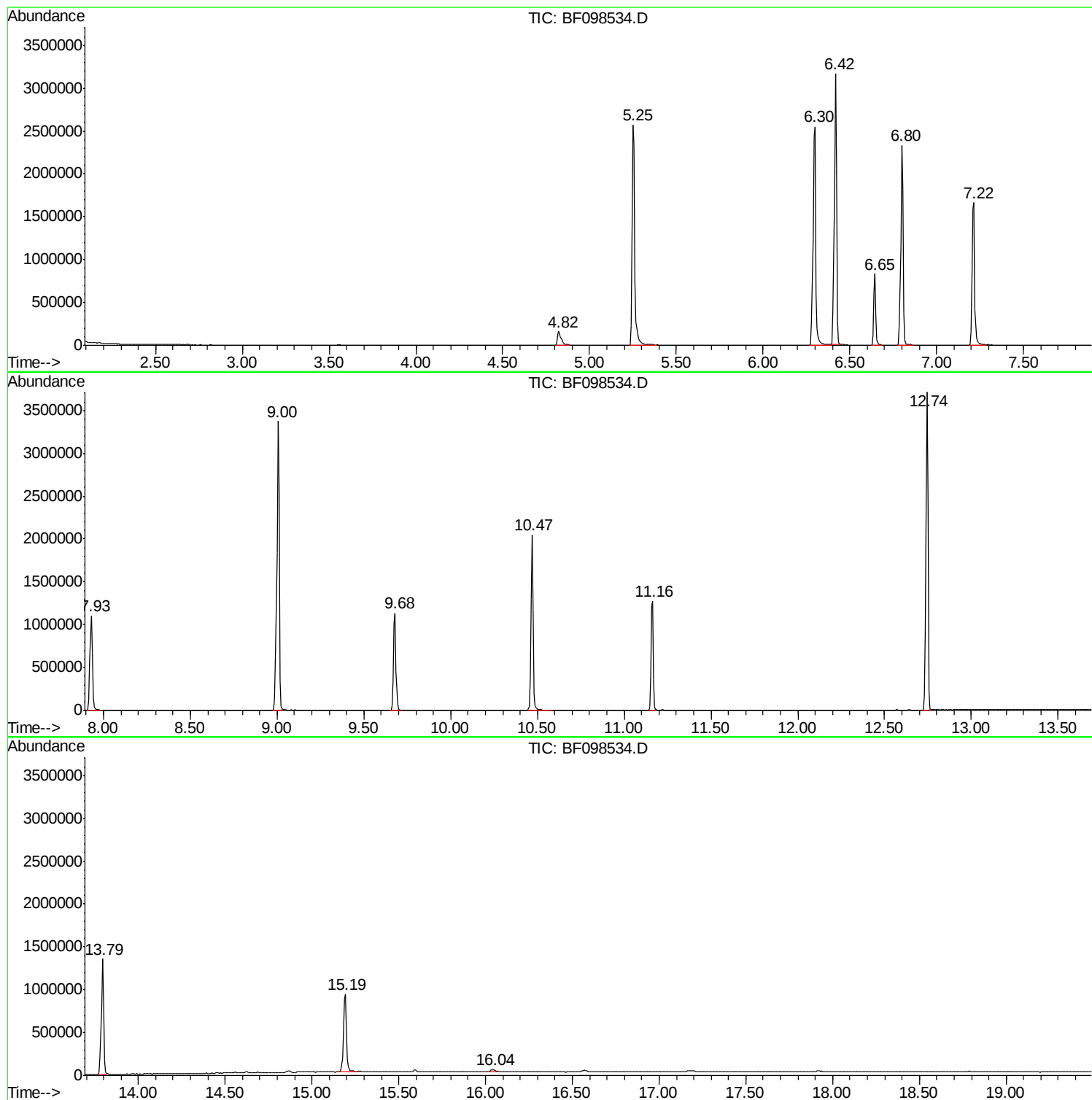
Sum of corrected areas: 27167966

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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF091117.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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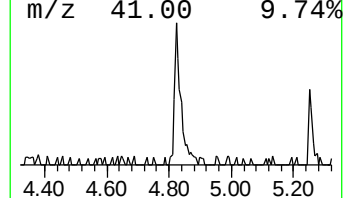
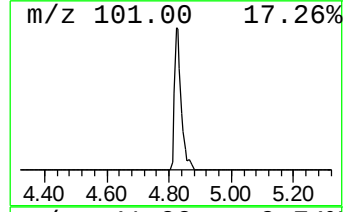
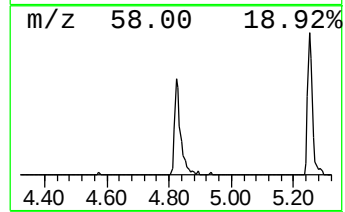
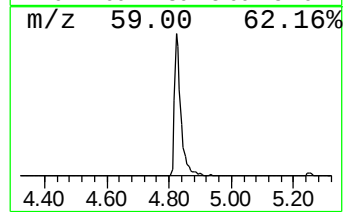
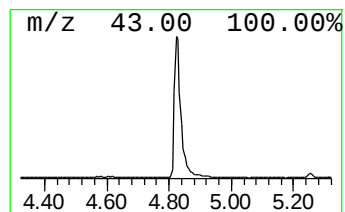
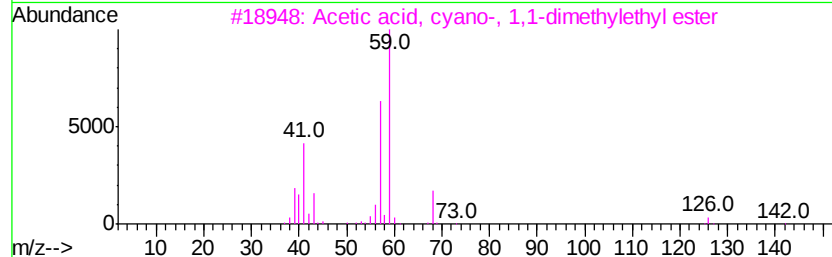
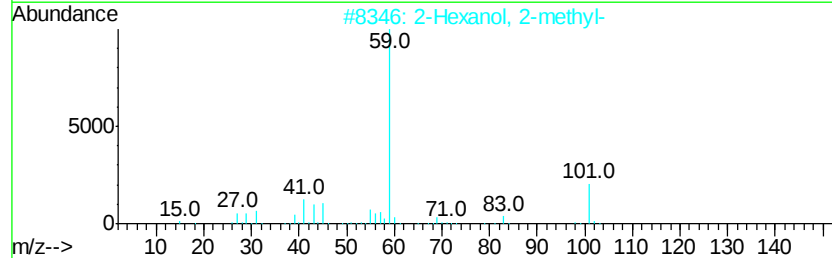
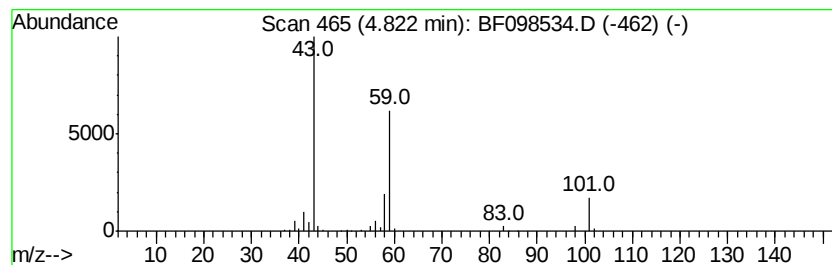
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.82	6.40 ng	227444	1,4-Dichlorobenzene-d4	6.65

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	59	
2	2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	28	
3	Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	23	
4	Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9	
5	5-Hexen-2-one	98	C6H10O	000109-49-9	9	



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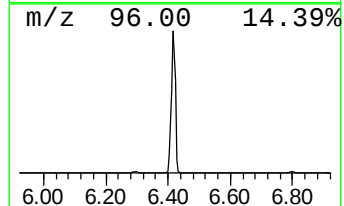
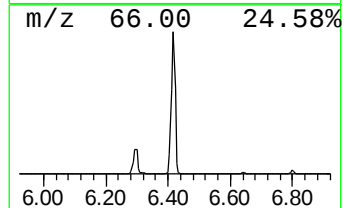
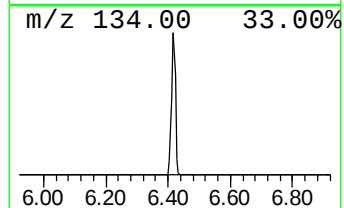
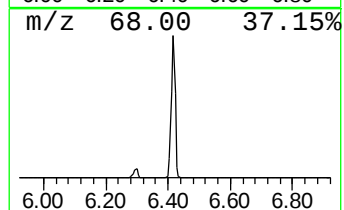
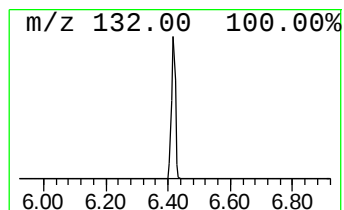
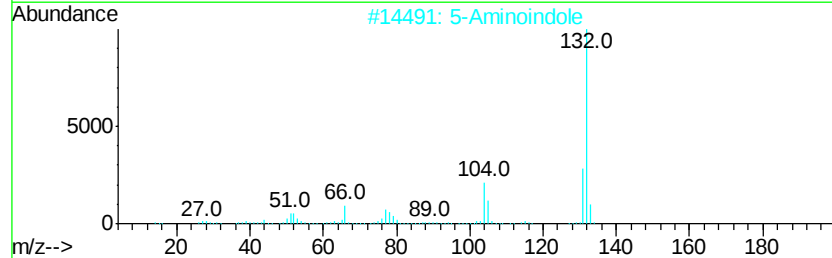
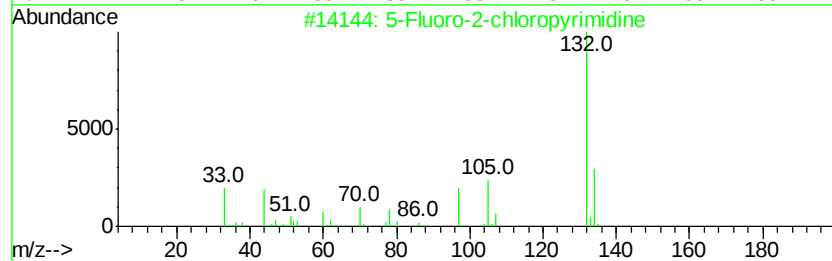
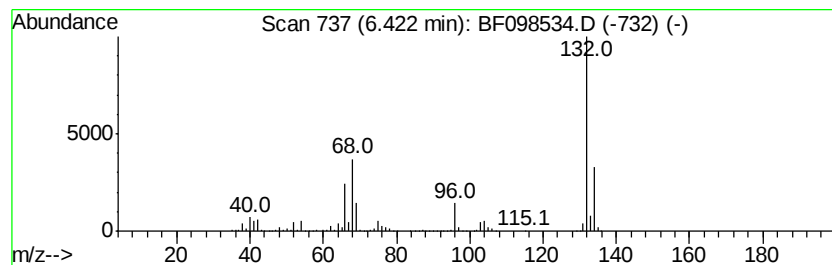
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Peak Number 2 unknown6.42 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.42	77.63 ng	2756920	1,4-Dichlorobenzene-d4	6.65

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	16
2			5-Aminoindole	132	C8H8N2	005192-03-0	12
3			Tranvlcyproline-propionyl	189	C12H15NO	1000123-86-3	10
4			Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	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...	4.82	6.4	ng	227444	1	6.65	710272	20.0
unknown6.42	6.42	77.6	ng	2756920	1	6.65	710272	20.0