

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF061518\
 Data File : BF106494.D
 Acq On : 15 Jun 2018 16:21
 Operator : JU/SJ
 Sample : J3467-12
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 C-MW-09-061218

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-BF061118.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	5.195	482	486	496	rVB	91596	98209	3.29%	0.463%
2	5.613	553	557	568	rVB	1393425	1416881	47.44%	6.682%
3	6.607	723	726	740	rBV	1016328	937694	31.39%	4.422%
4	6.742	745	749	753	rBV	2540274	2362697	79.10%	11.143%
5	6.966	783	787	790	rBV	943143	743643	24.90%	3.507%
6	7.125	809	814	817	rBV	2461110	2279928	76.33%	10.753%
7	7.531	877	883	886	rBV	2011894	1926228	64.49%	9.085%
8	8.248	1001	1005	1008	rBV	1084410	871780	29.19%	4.112%
9	9.325	1183	1188	1191	rBV	3138046	2986857	100.00%	14.087%
10	9.713	1249	1254	1259	rBV	150928	128550	4.30%	0.606%
11	9.919	1286	1289	1293	rBV	49917	39720	1.33%	0.187%
12	10.007	1300	1304	1310	rVB	979196	848408	28.40%	4.001%
13	10.419	1372	1374	1377	rVB	64433	48821	1.63%	0.230%
14	10.801	1434	1439	1442	rBV	1687090	1547010	51.79%	7.296%
15	10.895	1451	1455	1464	rVB	49296	51562	1.73%	0.243%
16	11.495	1554	1557	1561	rBV	788942	712794	23.86%	3.362%
17	13.083	1823	1827	1830	rBV	2768420	2521428	84.42%	11.892%
18	13.483	1891	1895	1898	rBV	51184	42347	1.42%	0.200%
19	13.971	1974	1978	1986	rBV3	33631	43592	1.46%	0.206%
20	14.154	2005	2009	2013	rBV	938077	842039	28.19%	3.971%
21	15.701	2267	2272	2278	rVB2	577188	753058	25.21%	3.552%

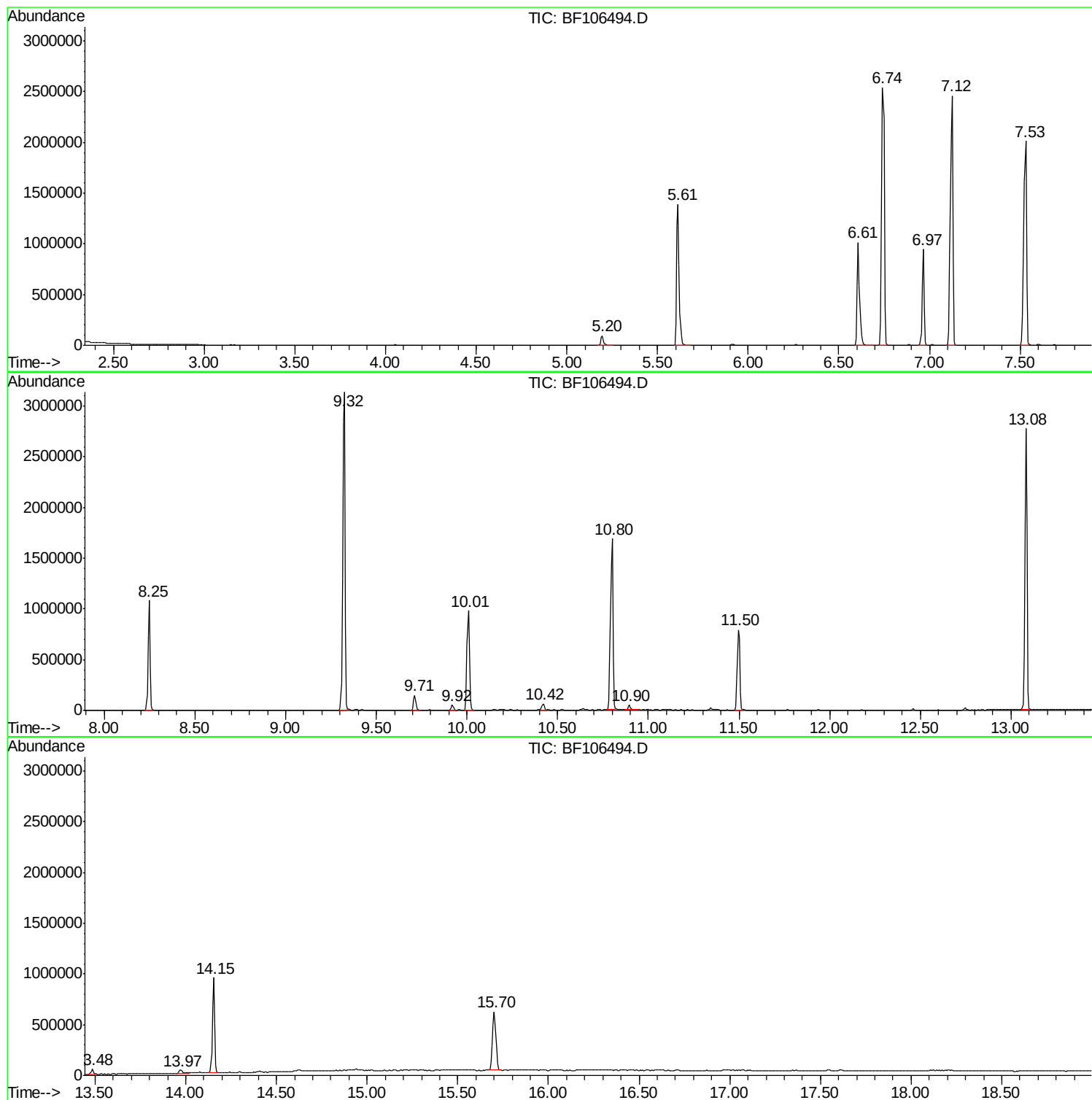
Sum of corrected areas: 21203246

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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF061118.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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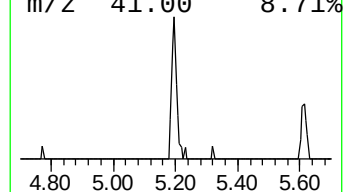
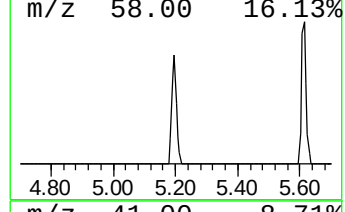
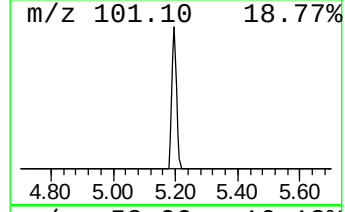
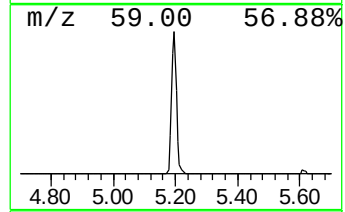
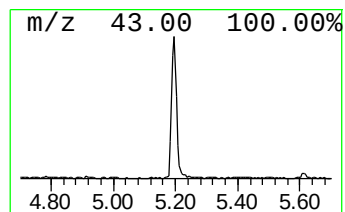
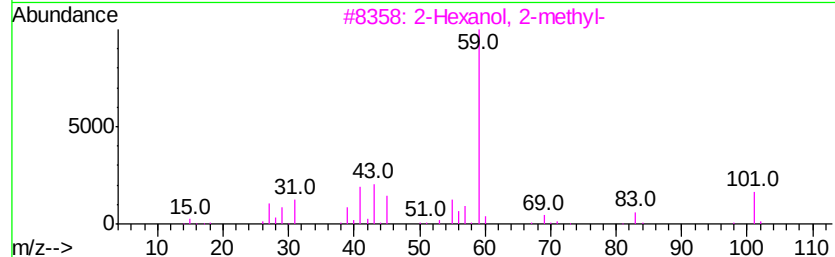
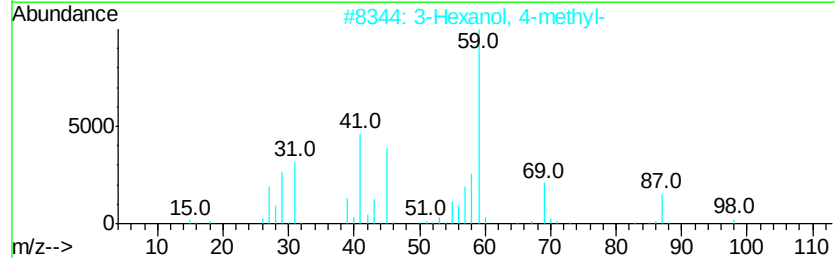
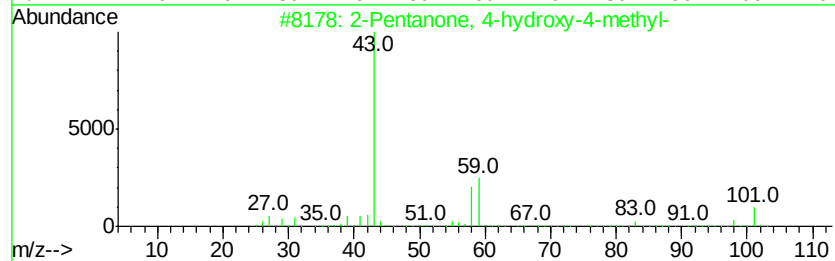
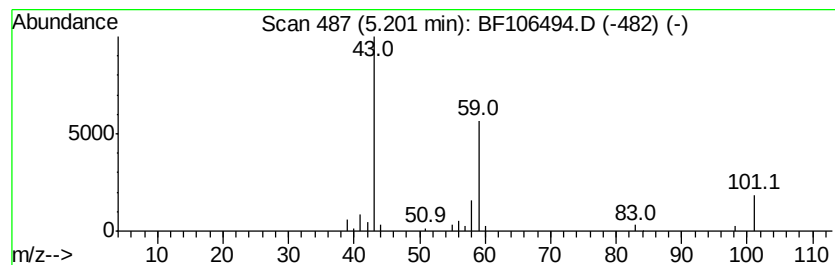
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF061118.M
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.20	2.64 ng	98209	1,4-Dichlorobenzene-d4	6.97

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	59
2			3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	33
3			2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	28
4			Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	25
5			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9



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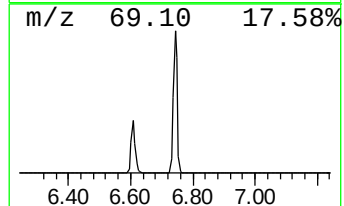
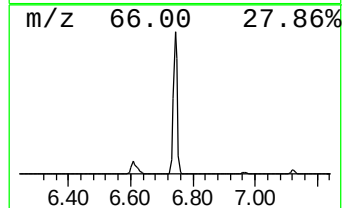
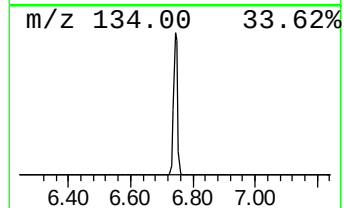
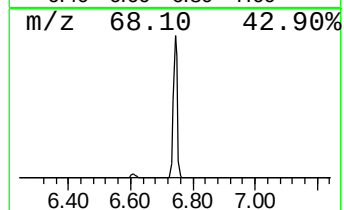
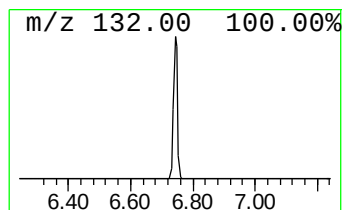
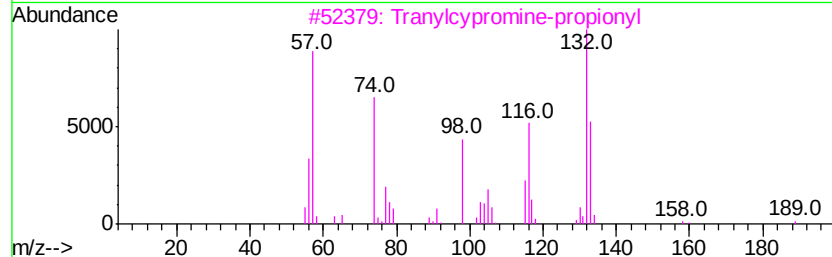
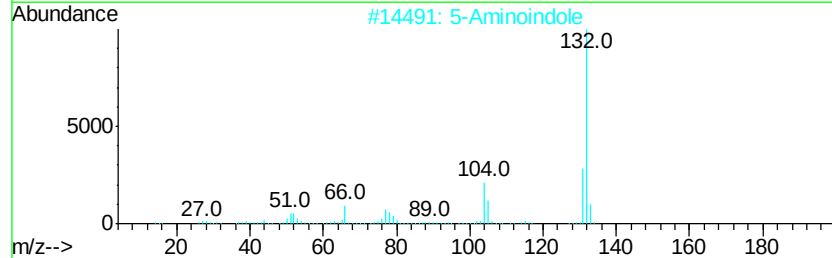
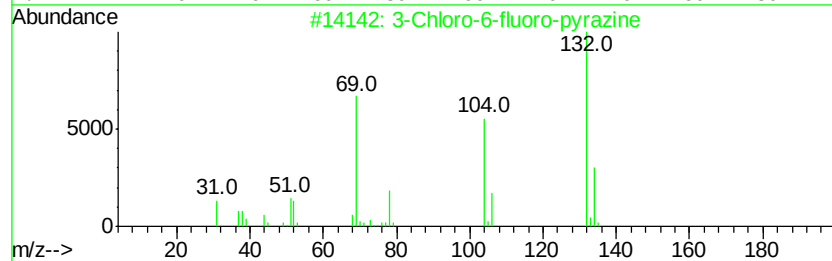
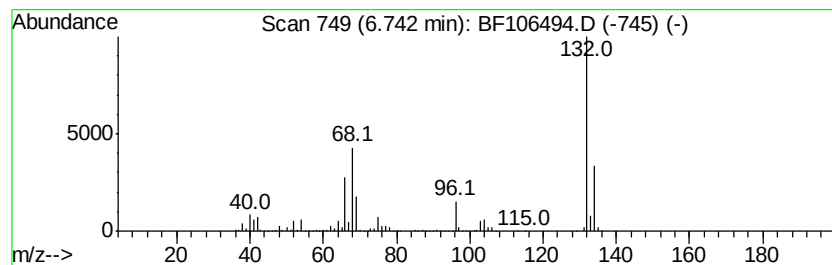
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Peak Number 2 unknown6.74 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.74	63.54 ng	2362700	1,4-Dichlorobenzene-d4	6.97

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			Tranvlcypropine-propionyl	189	C12H15NO	1000123-86-3	10
4			5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	9
5			1,3,4-Thiadiazole-2(3H)-thione, ...	132	C3H4N2S2	029490-19-5	9



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TIC Top Hit name	RT	EstConc	Units	Response	---Internal Standard---			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.20	2.6	ng	98209	1	6.97	743643	20.0
unknown6.74	6.74	63.5	ng	2362700	1	6.97	743643	20.0