

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF082818\
 Data File : BF108581.D
 Acq On : 29 Aug 2018 3:22
 Operator : JU/SJ
 Sample : J4656-04
 Misc :
 ALS Vial : 28 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 082418-SIDI-F-U-B

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-BF080818.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	5.237	490	493	505	rBV	57808	84592	2.12%	0.390%
2	5.637	558	561	589	rBV	389554	899913	22.53%	4.147%
3	6.631	727	730	750	rBV	292289	649210	16.25%	2.992%
4	6.772	750	754	779	rVB	1612283	1935912	48.47%	8.921%
5	7.001	789	793	815	rVB	521684	679307	17.01%	3.130%
6	7.154	815	819	851	rBV	2605297	2843620	71.20%	13.104%
7	7.566	884	889	912	rBV	2042488	2332038	58.39%	10.747%
8	8.284	1007	1011	1026	rBV	661062	792496	19.84%	3.652%
9	9.354	1189	1193	1203	rBV	4126524	3994106	100.00%	18.406%
10	9.748	1256	1260	1271	rBV	133093	156798	3.93%	0.723%
11	10.042	1307	1310	1321	rVB	856662	806633	20.20%	3.717%
12	10.695	1418	1421	1428	rBV	80735	79512	1.99%	0.366%
13	10.836	1441	1445	1458	rBV	1290332	1335111	33.43%	6.152%
14	11.536	1561	1564	1573	rBV	694685	702449	17.59%	3.237%
15	13.124	1830	1834	1837	rBV	3354971	3045803	76.26%	14.036%
16	14.195	2011	2016	2021	rBV	780972	774156	19.38%	3.567%
17	15.748	2275	2280	2289	rVB	435488	588727	14.74%	2.713%

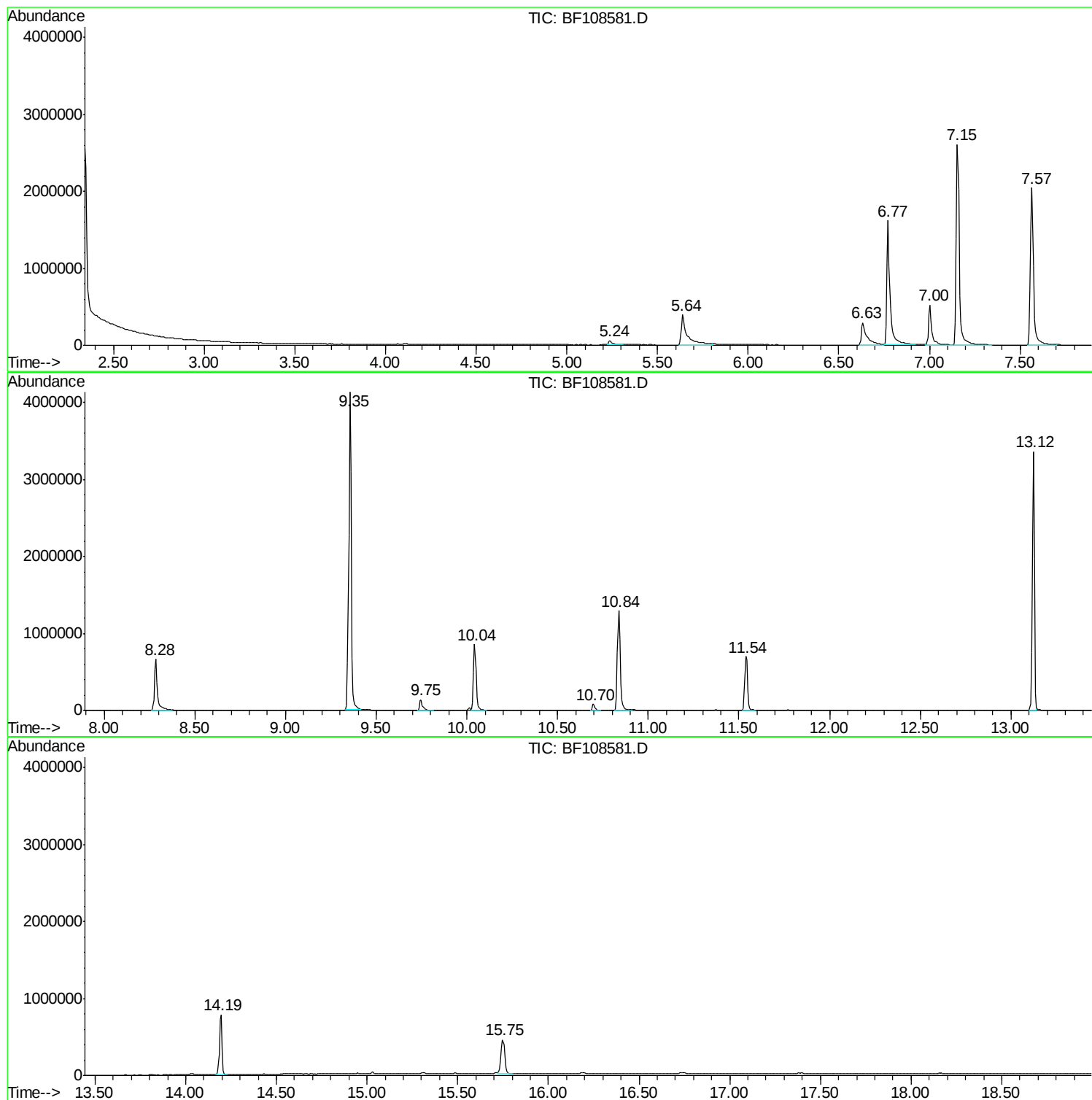
Sum of corrected areas: 21700383

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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF080818.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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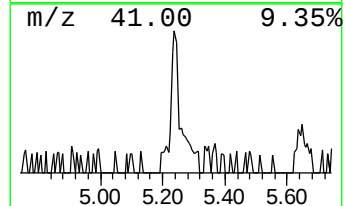
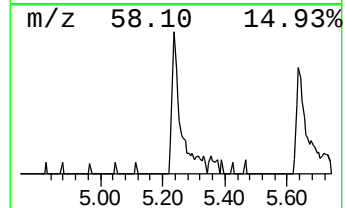
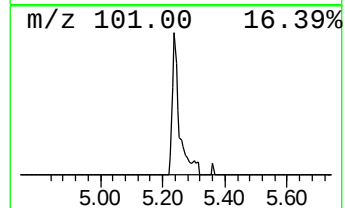
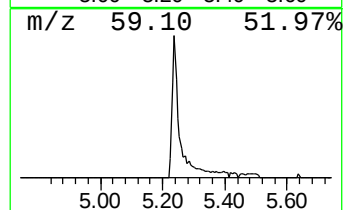
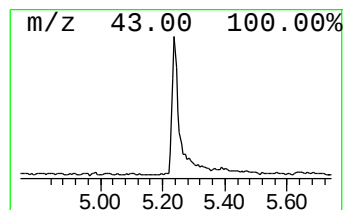
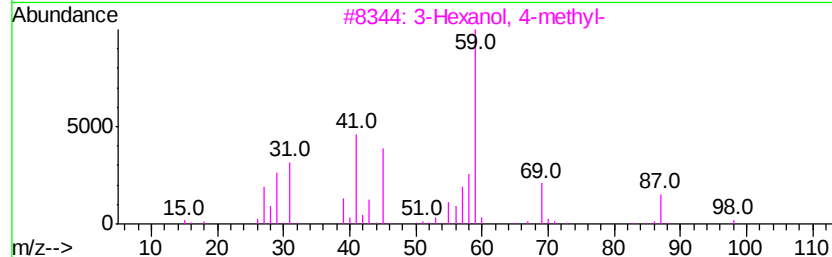
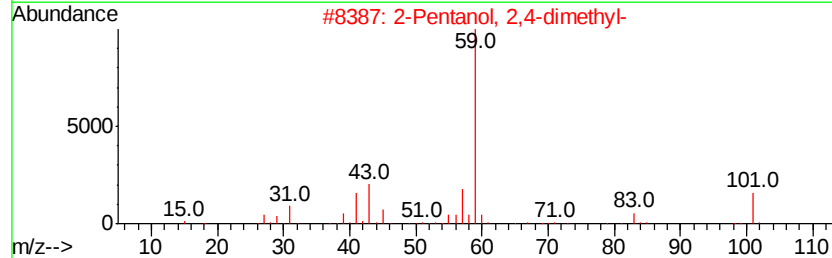
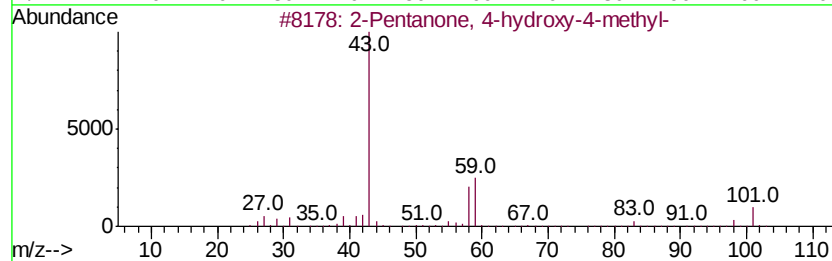
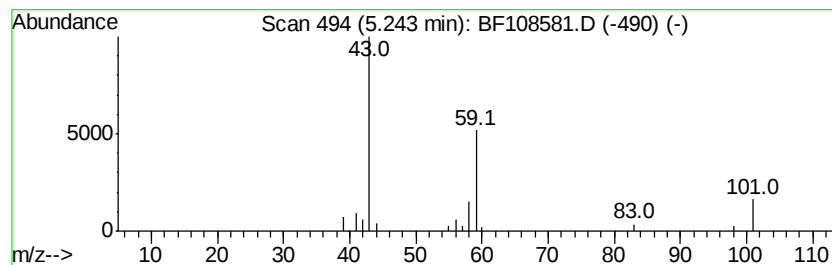
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF080818.M
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TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.24	2.49 ng	84592	1,4-Dichlorobenzene-d4	7.00

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	59
2		2-Pentanol, 2,4-dimethyl-	116	C7H16O	000625-06-9	40
3		3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	28
4		2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	28
5		Butyl aldoxime, 3-methyl-, syn-	101	C5H11NO	005780-40-5	23



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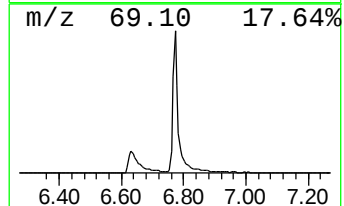
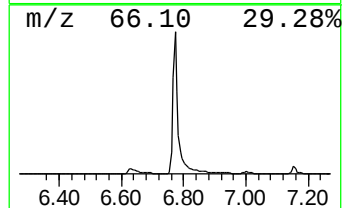
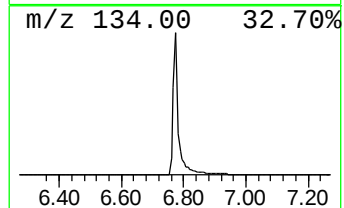
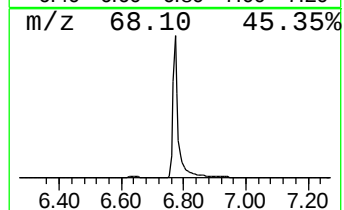
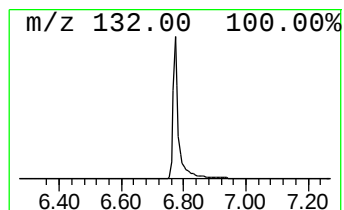
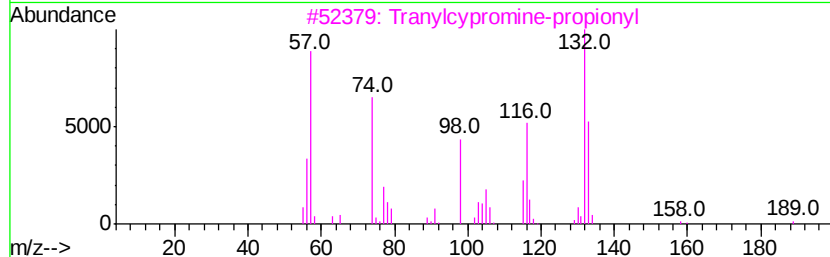
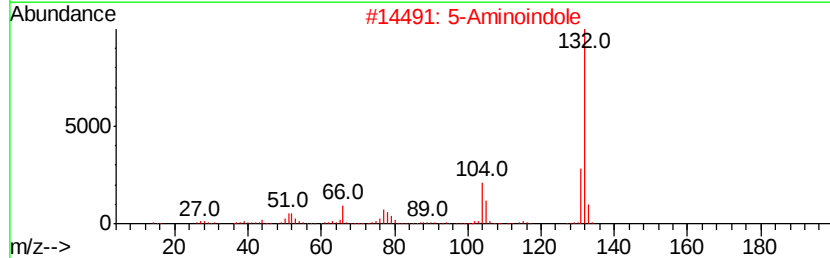
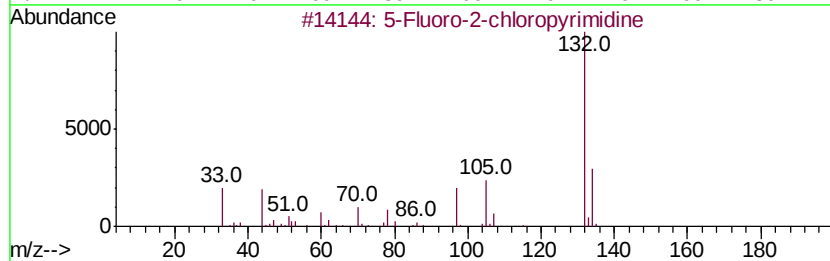
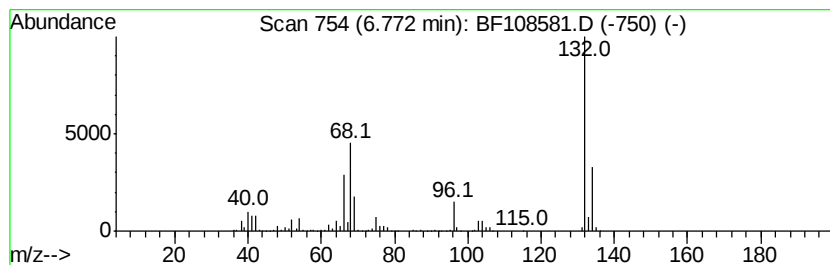
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Peak Number 2 unknown6.77 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.77	57.00 ng	1935910	1,4-Dichlorobenzene-d4	7.00

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	16
2		5-Aminoindole	132	C8H8N2	005192-03-0	12
3		Tranvlcypromine-propionyl	189	C12H15NO	1000123-86-3	10
4		(5-Methyl-2-pyridyl)acetonitrile	132	C8H8N2	1000241-93-9	10
5		1,2,4-Trifluorobenzene	132	C6H3F3	000367-23-7	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.24	2.5	ng	84592	1	7.00	679307	20.0
unknown6.77	6.77	57.0	ng	1935910	1	7.00	679307	20.0