

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF072518\
 Data File : BF107688.D
 Acq On : 26 Jul 2018 6:23
 Operator : JU/SJ
 Sample : J4133-05
 Misc :
 ALS Vial : 39 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 072318-DBRI-F-U-M

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-BF072018.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	5.596	550	554	579	rBV	1006466	1556365	26.52%	4.766%
2	6.595	721	724	744	rBV	631673	1200225	20.45%	3.676%
3	6.737	744	748	772	rVB	2741516	3132151	53.36%	9.592%
4	6.966	784	787	794	rBV	851141	1008126	17.18%	3.087%
5	7.125	809	814	841	rBV	4076892	4526466	77.12%	13.862%
6	7.537	879	884	908	rBV	2678925	3580479	61.00%	10.965%
7	8.248	1001	1005	1022	rBV	1151898	1339537	22.82%	4.102%
8	9.325	1183	1188	1219	rBV	5297594	5869578	100.00%	17.976%
9	9.713	1251	1254	1260	rBV	68327	71219	1.21%	0.218%
10	10.007	1301	1304	1325	rVB	1269283	1255785	21.39%	3.846%
11	10.666	1413	1416	1424	rBV	284282	237936	4.05%	0.729%
12	10.801	1435	1439	1453	rBV	1837347	2065701	35.19%	6.326%
13	11.501	1554	1558	1567	rBV	1032016	1106119	18.84%	3.388%
14	13.089	1823	1828	1846	rBV	3380035	3623023	61.73%	11.096%
15	13.977	1974	1979	1988	rBV4	29816	64191	1.09%	0.197%
16	14.154	2005	2009	2019	rBV	900569	974057	16.60%	2.983%
17	15.265	2194	2198	2201	rVB	58616	66869	1.14%	0.205%
18	15.689	2263	2270	2287	rVB	533190	901268	15.35%	2.760%
19	16.130	2341	2345	2351	rBV2	50431	73741	1.26%	0.226%

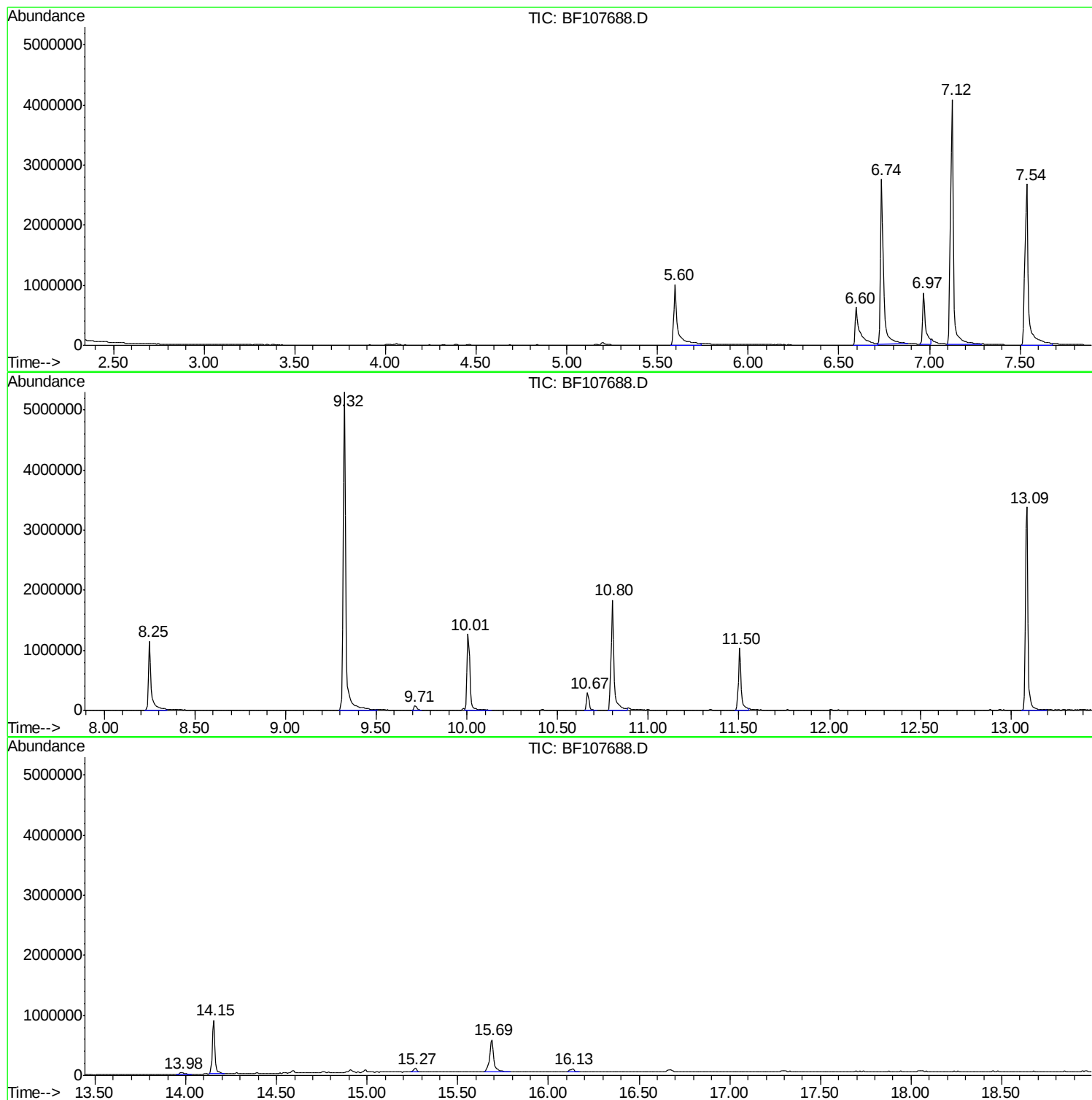
Sum of corrected areas: 32652836

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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF072018.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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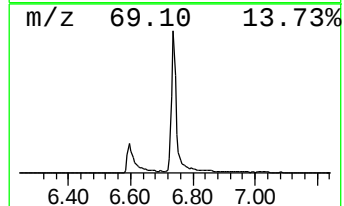
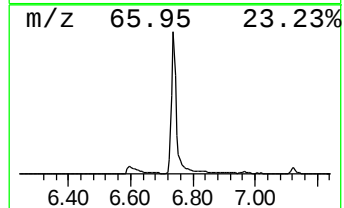
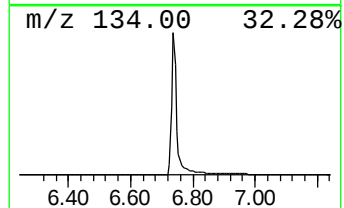
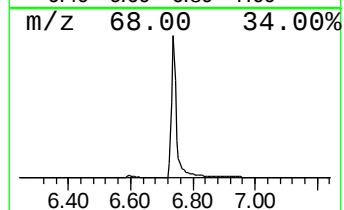
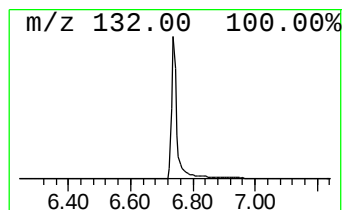
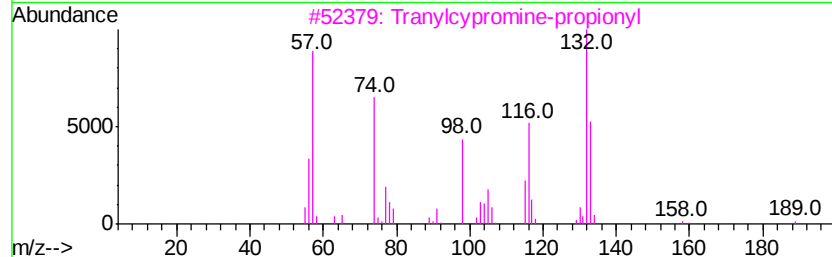
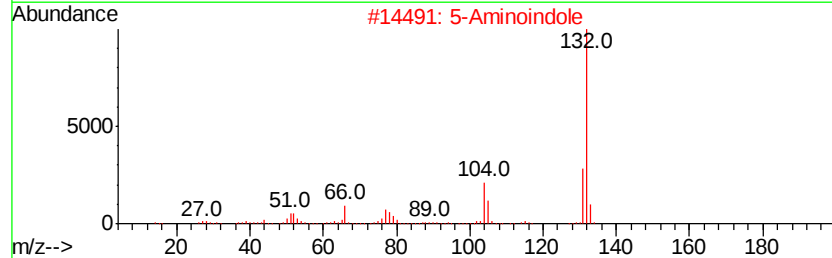
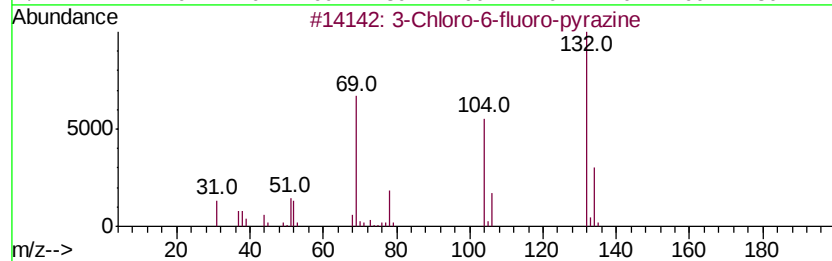
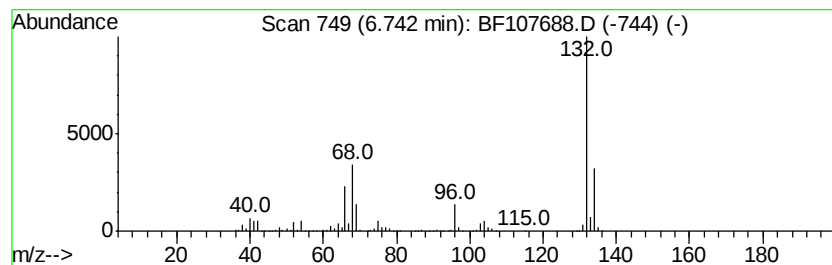
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TIC Integration Parameters: LSCINT.P

Peak Number 1 unknown6.74 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.74	62.14 ng	3132150	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			Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9
5			2H-Cyclopenta[d]pyridazine, 2-me...	132	C8H8N2	022291-85-6	9



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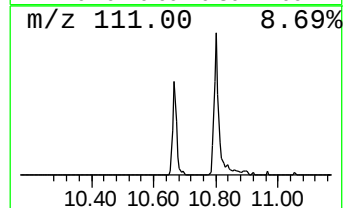
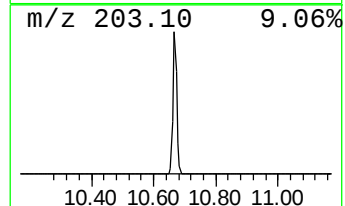
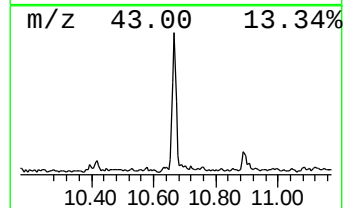
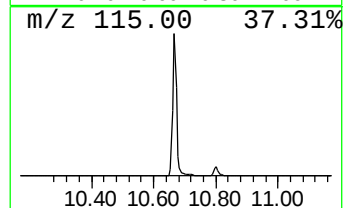
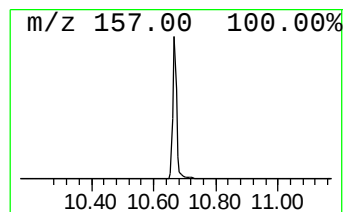
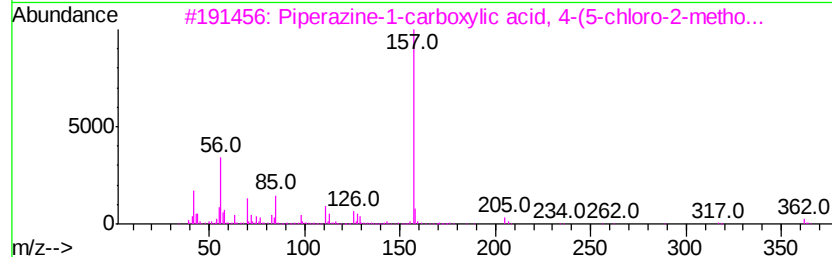
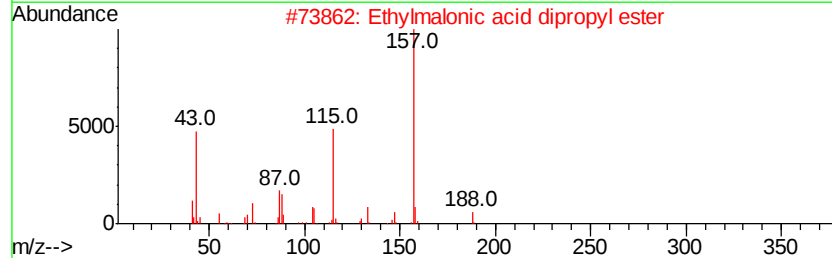
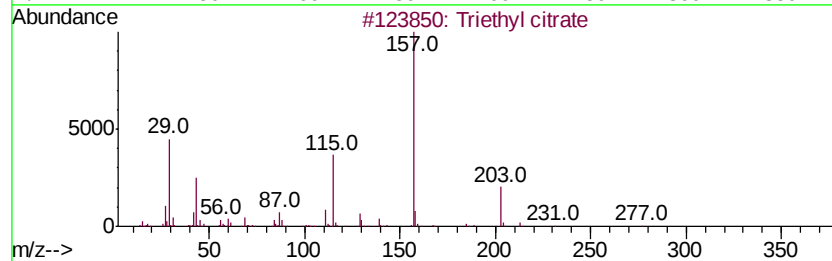
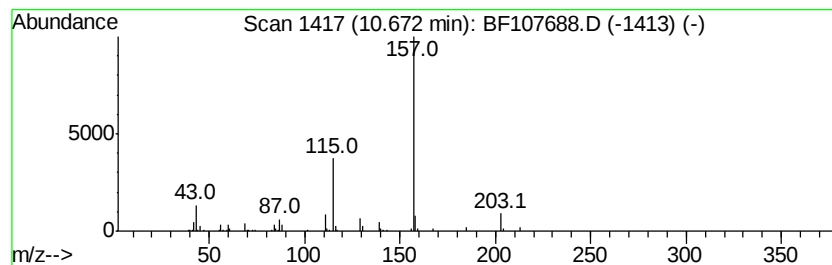
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Peak Number 3 Triethyl citrate Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.67	3.79 ng	237936	Acenaphthene-d10	10.01
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Triethyl citrate	276 C12H20O7	000077-93-0 91
2		Ethylmalonic acid dipropyl ester	216 C11H20O4	001113-91-3 50
3		Piperazine-1-carboxylic acid, 4-...	362 C14H19ClN2O5S	1000303-80-2 42
4		Quinoline, 4,8-dimethyl-	157 C11H11N	013362-80-6 40
5		3-Chloro2-fluorobenzoic acid, 3-...	370 C21H32ClF02	1000338-64-6 36



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TIC Top Hit name	RT	EstConc	Units	Response	---Internal Standard---			
					#	RT	Resp	Conc
unknown6.74	6.74	62.1	ng	3132150	1	6.97	1008130	20.0
Triethyl citrate	10.67	3.8	ng	237936	3	10.01	1255790	20.0