

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF062618\
 Data File : BF106843.D
 Acq On : 27 Jun 2018 1:10
 Operator : JU/SJ
 Sample : J3639-02
 Misc :
 ALS Vial : 29 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 062018-DBRI-E-D-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-BF061918.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	4.090	294	298	303	rBB	18760	22243	1.87%	0.328%
2	5.190	482	485	490	rBV	16301	16825	1.42%	0.248%
3	5.601	551	555	571	rBV	325580	299559	25.21%	4.412%
4	6.601	721	725	740	rBB2	240006	262585	22.10%	3.867%
5	6.737	744	748	754	rBV	569157	533313	44.88%	7.854%
6	6.960	782	786	789	rBV	272640	234680	19.75%	3.456%
7	7.113	808	812	816	rBV	833277	799174	67.26%	11.770%
8	7.525	877	882	885	rBV	626695	577317	48.59%	8.502%
9	8.242	1000	1004	1019	rBV	316643	268109	22.56%	3.949%
10	9.313	1182	1186	1196	rBV	1160007	1082410	91.10%	15.941%
11	9.707	1250	1253	1259	rBB	28113	23632	1.99%	0.348%
12	10.001	1300	1303	1312	rVB	286210	264947	22.30%	3.902%
13	10.407	1369	1372	1375	rVB	20878	15656	1.32%	0.231%
14	10.660	1412	1415	1418	rBV	28866	23419	1.97%	0.345%
15	10.795	1434	1438	1445	rBV	334517	333740	28.09%	4.915%
16	10.883	1450	1453	1458	rBV	12326	14012	1.18%	0.206%
17	11.489	1553	1556	1564	rVB	287296	273695	23.03%	4.031%
18	13.077	1821	1826	1829	rBV	1260600	1188182	100.00%	17.499%
19	14.142	2003	2007	2012	rBV	328650	307230	25.86%	4.525%
20	14.989	2147	2151	2155	rVB	39646	42562	3.58%	0.627%
21	15.683	2263	2269	2278	rVB	155422	206679	17.39%	3.044%

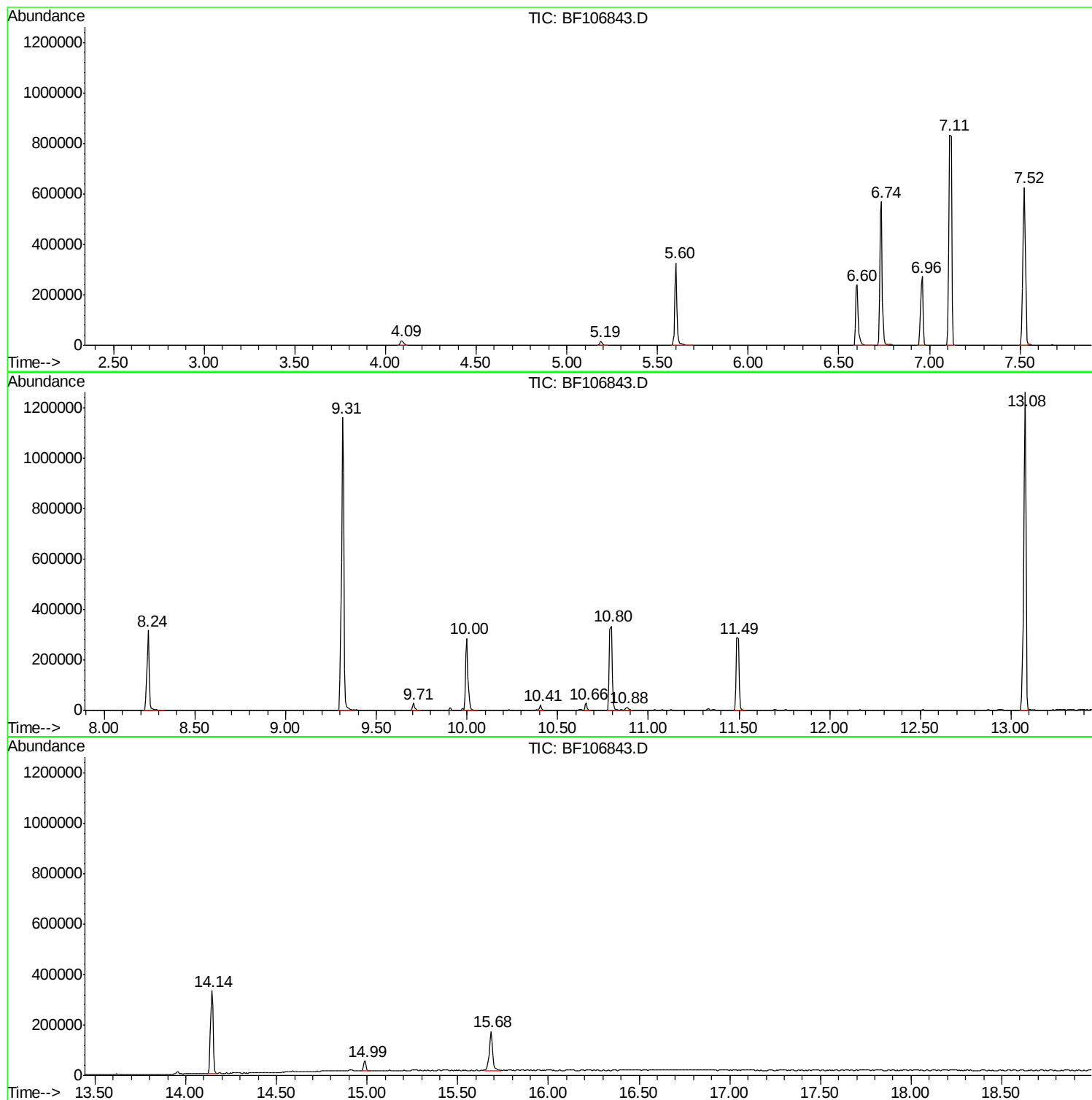
Sum of corrected areas: 6789969

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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF061918.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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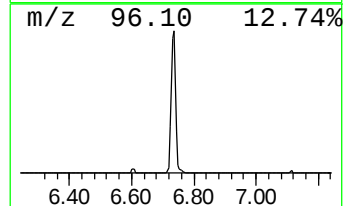
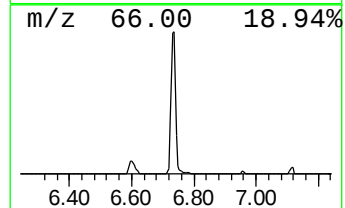
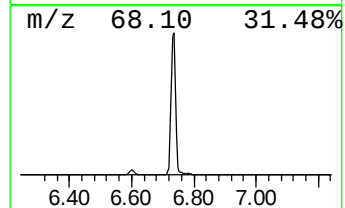
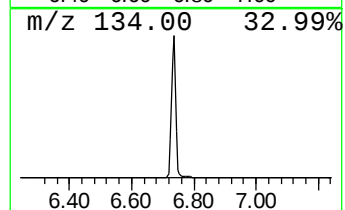
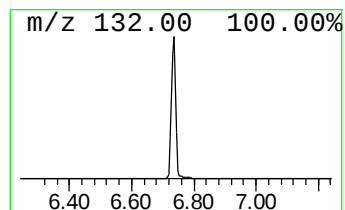
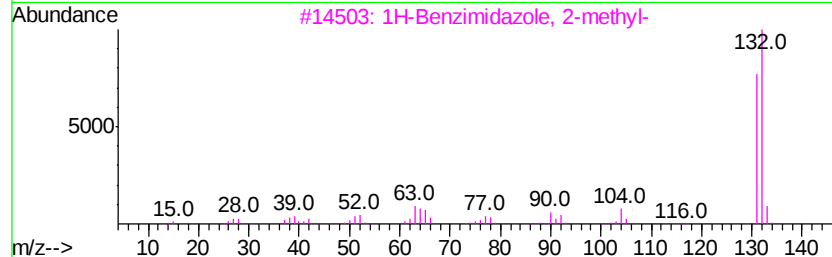
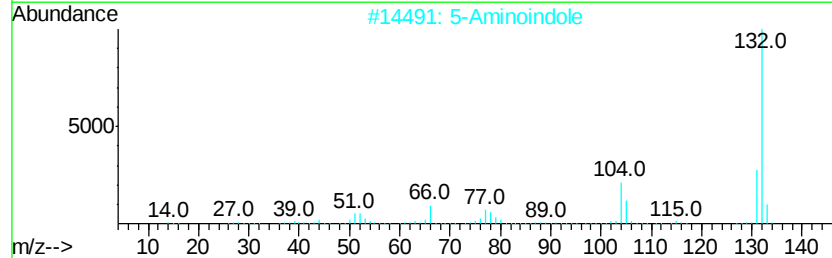
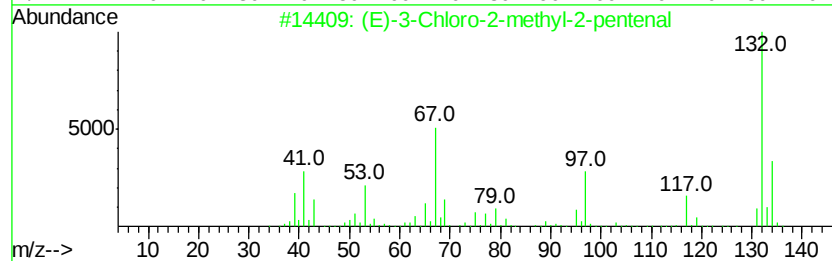
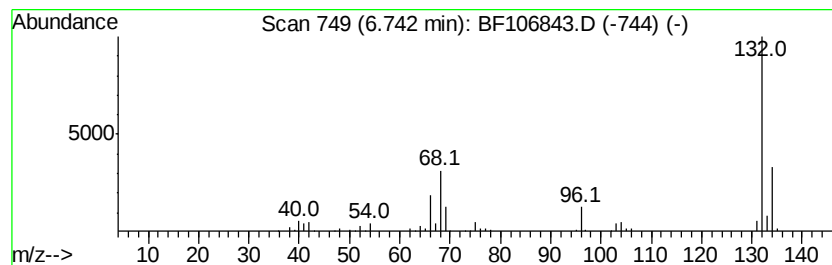
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF061918.M
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 unknown6.74 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.74	45.45 ng	533313	1,4-Dichlorobenzene-d4	6.96

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	17
2		5-Aminoindole	132	C8H8N2	005192-03-0	12
3		1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	10
4		Xenon	132	Xe	007440-63-3	9
5		4-Chloro-2-fluoro-pyrimidine	132	C4H2ClFN2	051422-00-5	9



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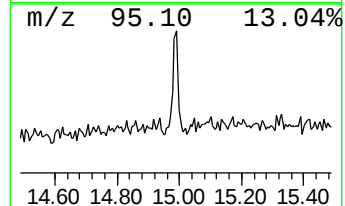
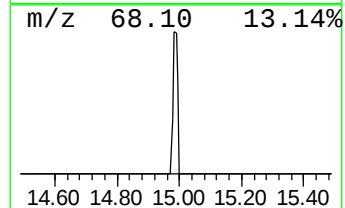
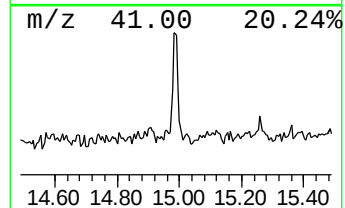
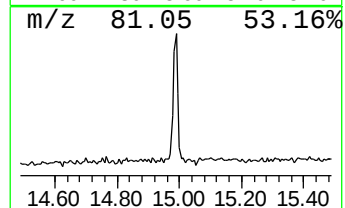
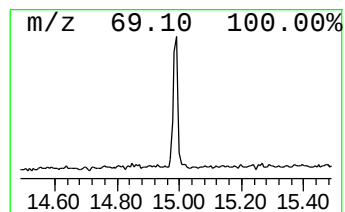
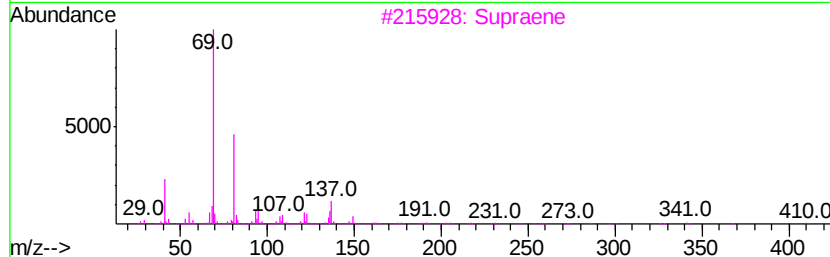
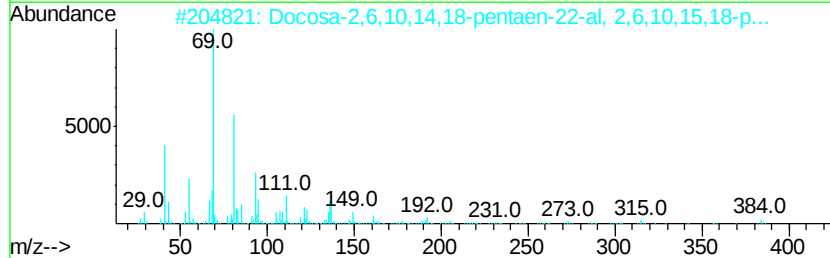
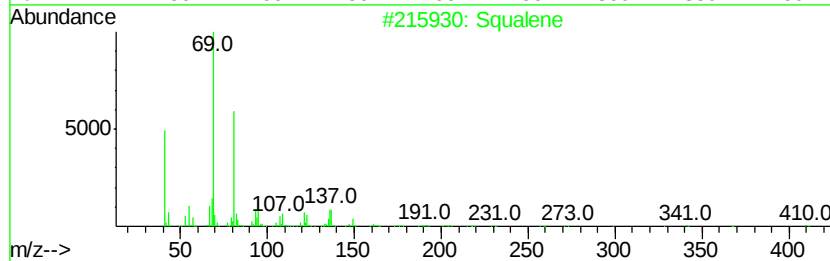
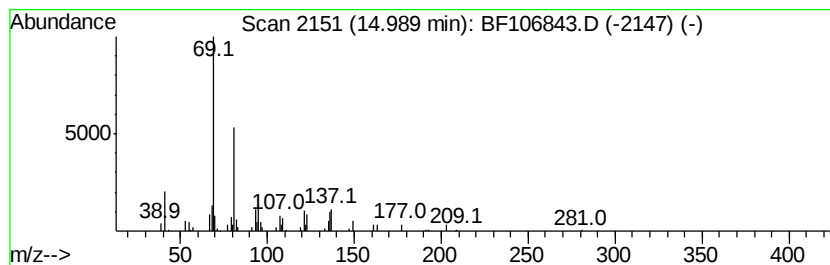
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Peak Number 3 Squalene Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.99	4.12 ng	42562	Perylene-d12	15.68

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Squalene	410	C30H50	000111-02-4	91
2		Docosa-2,6,10,14,18-pentaen-22-a...	384	C27H44O	1000163-04-7	91
3		Supraene	410	C30H50	007683-64-9	91
4		trans-Farnesol	222	C15H26O	000106-28-5	72
5		5,9,13-Pentadecatrien-2-one, 6,1...	262	C18H30O	001117-52-8	59



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
unknown6.74	6.74	45.5	ng	533313	1	6.96	234680	20.0
Squalene	14.99	4.1	ng	42562	6	15.68	206679	20.0