

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM031219\
 Data File : BM019138.D
 Acq On : 13 Mar 2019 08:28
 Operator : JU/SJ
 Sample : K1824-11
 Misc :
 ALS Vial : 30 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 TP-29

Integration Parameters: rteint.p
 Integrator: RTE
 Smoothing : ON
 Sampling : 1
 Start Thrs: 0.2
 Stop Thrs : 0

Filtering: 5
 Min Area: 3 % of largest Peak
 Max Peaks: 100
 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
 Peak separation: 5

Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\8270-BM031119.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	3.893	132	137	155	rVB	4930330	7076650	42.91%	7.801%
2	5.263	363	370	387	rBV	3875538	5954269	36.11%	6.564%
3	6.845	631	639	659	rBV	4362598	7278982	44.14%	8.024%
4	7.198	691	699	716	rBV	4347970	6849426	41.53%	7.551%
5	7.657	771	777	785	rBV	752250	1191604	7.23%	1.314%
6	7.975	823	831	846	rBV	2862200	4581634	27.78%	5.051%
7	8.828	968	976	1000	rBV	2203551	3892009	23.60%	4.291%
8	10.445	1244	1251	1272	rBV	940220	1676781	10.17%	1.848%
9	12.927	1666	1673	1697	rBV	6559079	9771593	59.26%	10.772%
10	13.780	1811	1818	1832	rBV	382265	626418	3.80%	0.691%
11	14.315	1901	1909	1922	rBV2	1723335	2639030	16.00%	2.909%
12	15.815	2157	2164	2182	rBV	6616362	9467530	57.41%	10.437%
13	17.068	2370	2377	2394	rBV	2367491	3452209	20.93%	3.806%
14	17.956	2523	2528	2539	rBV	177711	311965	1.89%	0.344%
15	19.709	2819	2826	2838	rBV	13026736	16490739	100.00%	18.179%
16	20.368	2935	2938	2942	rBV	137080	167454	1.02%	0.185%
17	20.962	3035	3039	3047	rBV	730999	873621	5.30%	0.963%
18	21.268	3085	3091	3101	rBV	2872222	3957906	24.00%	4.363%
19	21.403	3110	3114	3119	rBV	159158	201697	1.22%	0.222%
20	21.833	3184	3187	3194	rBV	171733	258733	1.57%	0.285%
21	22.291	3262	3265	3269	rVB	201922	245228	1.49%	0.270%
22	22.791	3347	3350	3356	rVB	201324	260433	1.58%	0.287%
23	23.356	3443	3446	3452	rVB	196063	282175	1.71%	0.311%
24	23.503	3465	3471	3481	rVB2	1616870	3203693	19.43%	3.532%

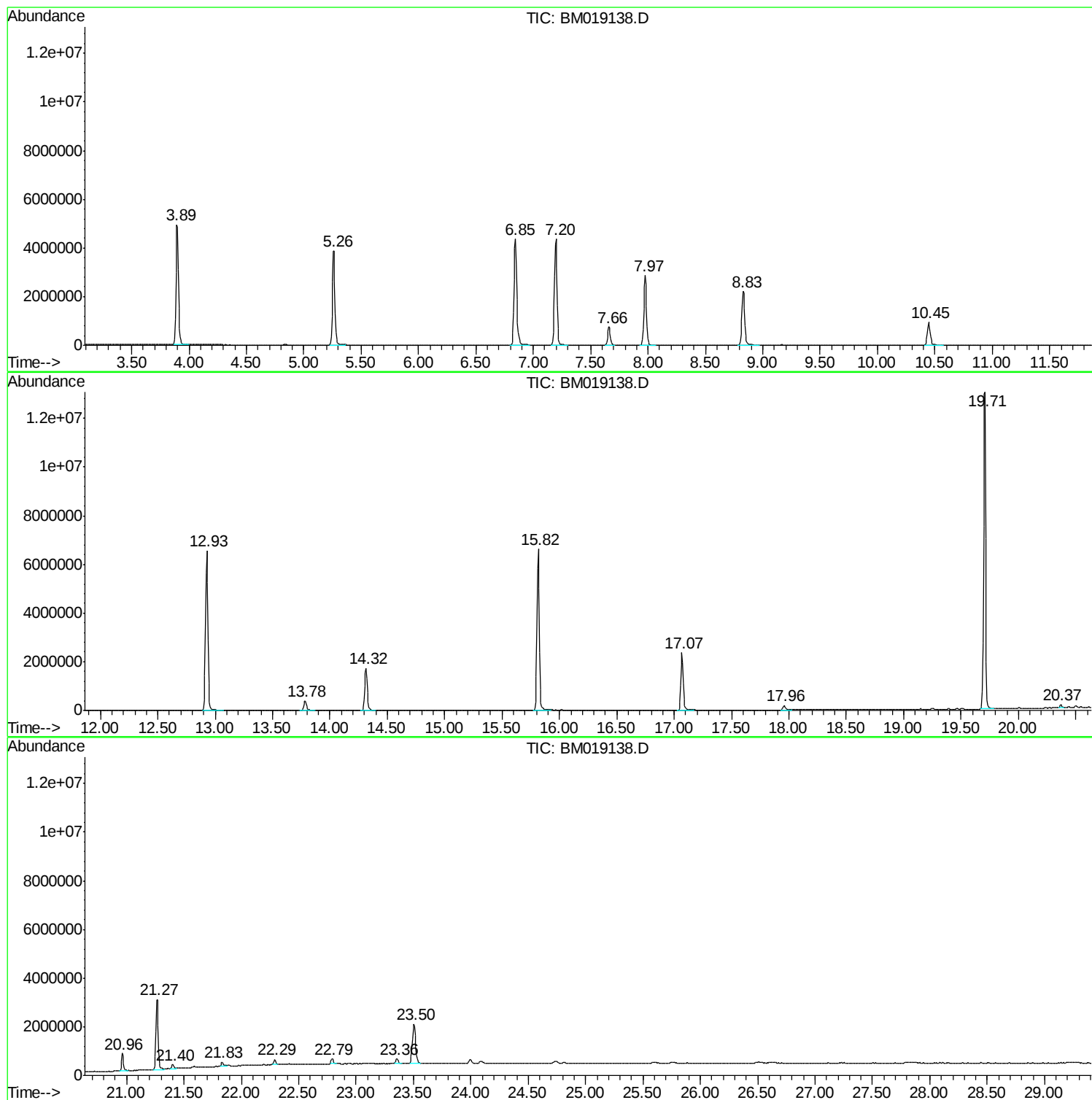
Sum of corrected areas: 90711779

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM031219\
Data File : BM019138.D
Acq On : 13 Mar 2019 08:28
Operator : JU/SJ
Sample : K1824-11
Misc :
ALS Vial : 30 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
TP-29

Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM031119.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM031219\
Data File : BM019138.D
Acq On : 13 Mar 2019 08:28
Operator : JU/SJ
Sample : K1824-11
Misc :
ALS Vial : 30 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
TP-29

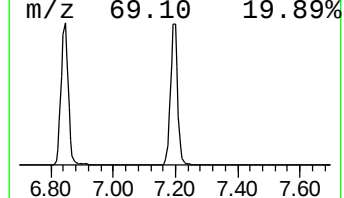
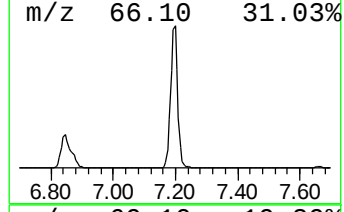
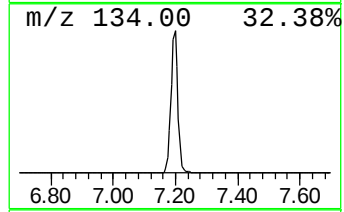
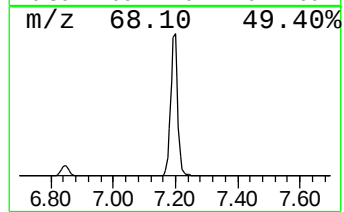
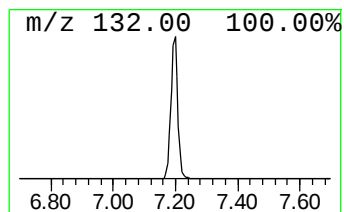
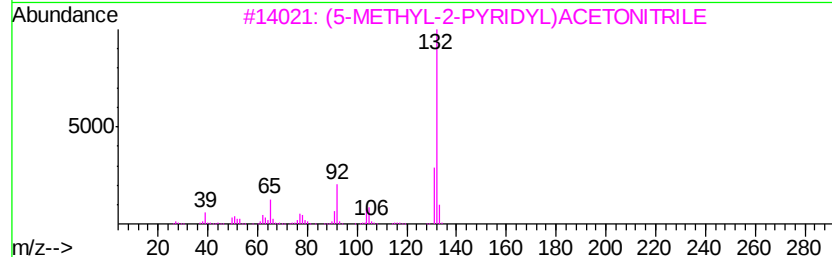
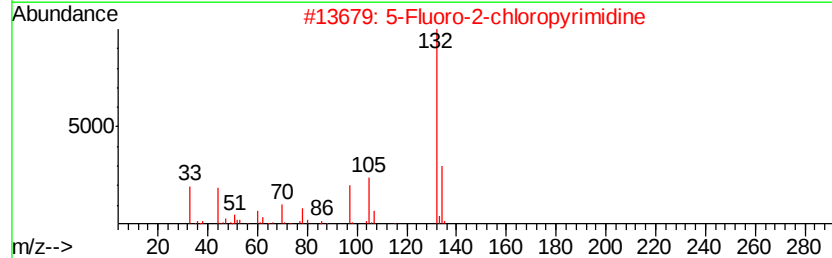
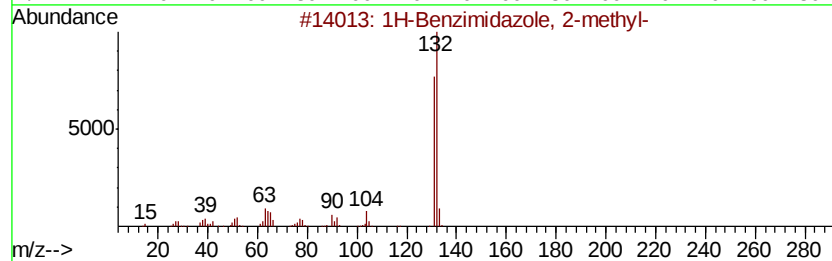
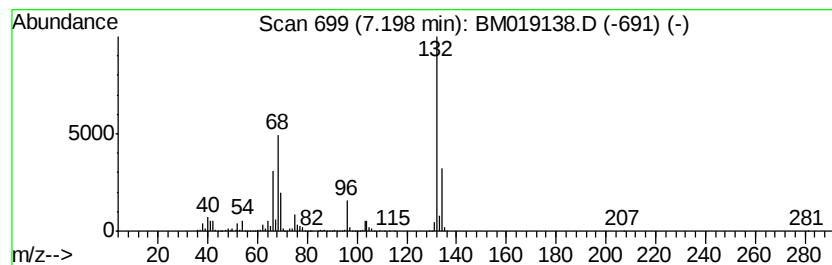
Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM031119.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 unknown7.20 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.20	114.96 ng	6849430	1,4-Dichlorobenzene-d4	7.66

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	22
2		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	12
3		(5-METHYL-2-PYRIDYL)ACETONITRILE	132	C8H8N2	1000241-93-9	10
4		5-Aminoindole	132	C8H8N2	005192-03-0	10
5		Benzene, 1,2,4-trifluoro-	132	C6H3F3	000367-23-7	9



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM031219\
 Data File : BM019138.D
 Acq On : 13 Mar 2019 08:28
 Operator : JU/SJ
 Sample : K1824-11
 Misc :
 ALS Vial : 30 Sample Multiplier: 1

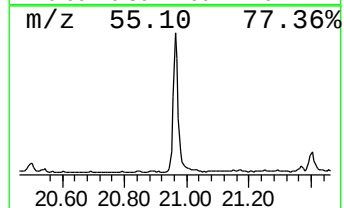
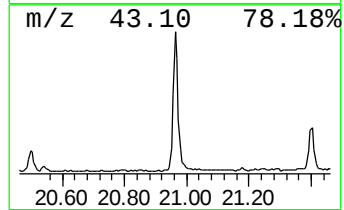
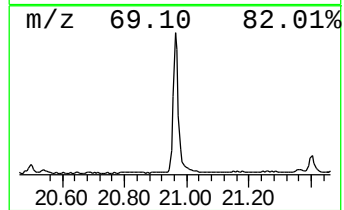
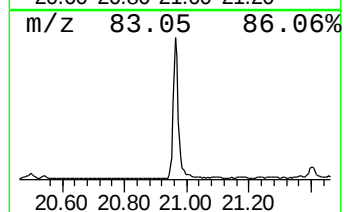
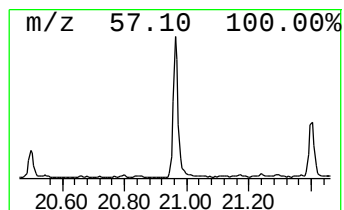
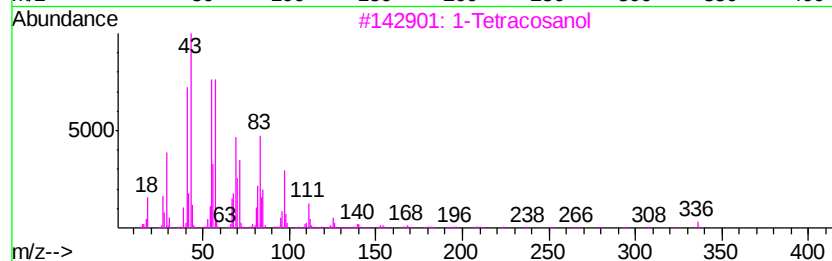
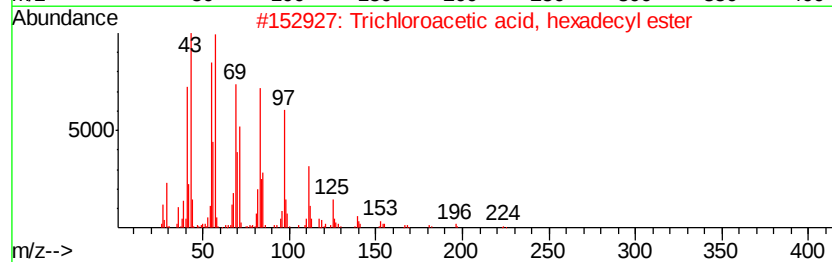
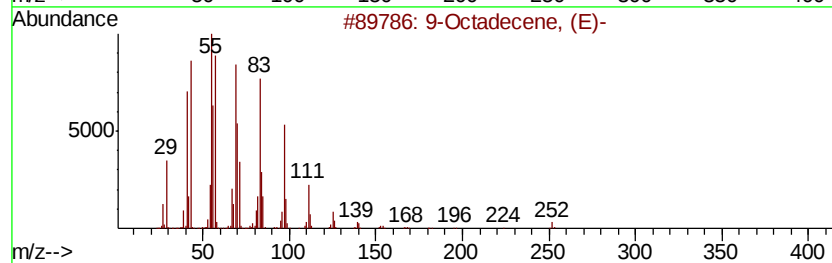
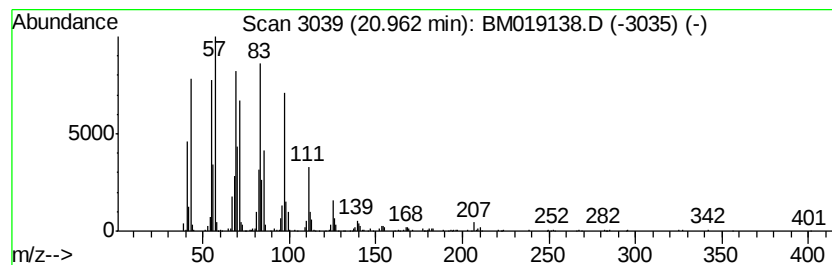
Instrument :
 BNA_M
 ClientSampleId :
 TP-29

Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM031119.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
 TIC Integration Parameters: LSCINT.P

 Peak Number 4 9-Octadecene, (E)- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.		
20.96	4.41 ng	873621	Chrysene-d12	21.27		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		9-Octadecene, (E)-	252	C18H36	007206-25-9	91
2		Trichloroacetic acid, hexadecyl ...	386	C18H33Cl3O2	074339-54-1	91
3		1-Tetracosanol	354	C24H50O	000506-51-4	91
4		Trichloroacetic acid, pentadecyl...	372	C17H31Cl3O2	074339-53-0	91
5		17-Pentatriacontene	491	C35H70	006971-40-0	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM031219\
Data File : BM019138.D
Acq On : 13 Mar 2019 08:28
Operator : JU/SJ
Sample : K1824-11
Misc :
ALS Vial : 30 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
TP-29

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\8270-BM031119.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
unknown7.20	7.20	115.0	ng	6849430	1	7.66	1191600	20.0
9-Octadecene, (E)-	20.96	4.4	ng	873621	5	21.27	3957910	20.0