

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM090618\
 Data File : BM016692.D
 Acq On : 06 Sep 2018 19:35
 Operator : SJ/JU
 Sample : J4798-08
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 090418-SIDI-E-U-B

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-BM082118.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	5.516	374	379	393	rBV2	305640	579795	4.18%	1.609%
2	7.146	651	656	671	rBV	199502	398854	2.88%	1.107%
3	7.493	709	715	734	rBV	809957	1436728	10.36%	3.987%
4	7.963	789	795	807	rBV2	259938	467841	3.37%	1.298%
5	8.275	841	848	864	rBV	1525969	2470618	17.82%	6.856%
6	9.146	989	996	1013	rBV	869315	1512004	10.91%	4.196%
7	10.763	1264	1271	1284	rBV	270857	507214	3.66%	1.408%
8	13.216	1680	1688	1701	rBV	3638286	4853693	35.01%	13.469%
9	14.069	1828	1833	1845	rBV	164093	238871	1.72%	0.663%
10	14.604	1917	1924	1931	rBV2	428349	701485	5.06%	1.947%
11	16.098	2172	2178	2193	rBV	2737147	3800327	27.41%	10.546%
12	17.345	2385	2390	2403	rVB	769924	1170699	8.44%	3.249%
13	18.204	2531	2536	2545	rBV3	149235	226445	1.63%	0.628%
14	19.939	2826	2831	2838	rBV	12183971	13862656	100.00%	38.470%
15	21.498	3091	3096	3105	rBV	1611763	2016549	14.55%	5.596%
16	23.756	3475	3480	3491	rVB	867422	1791366	12.92%	4.971%

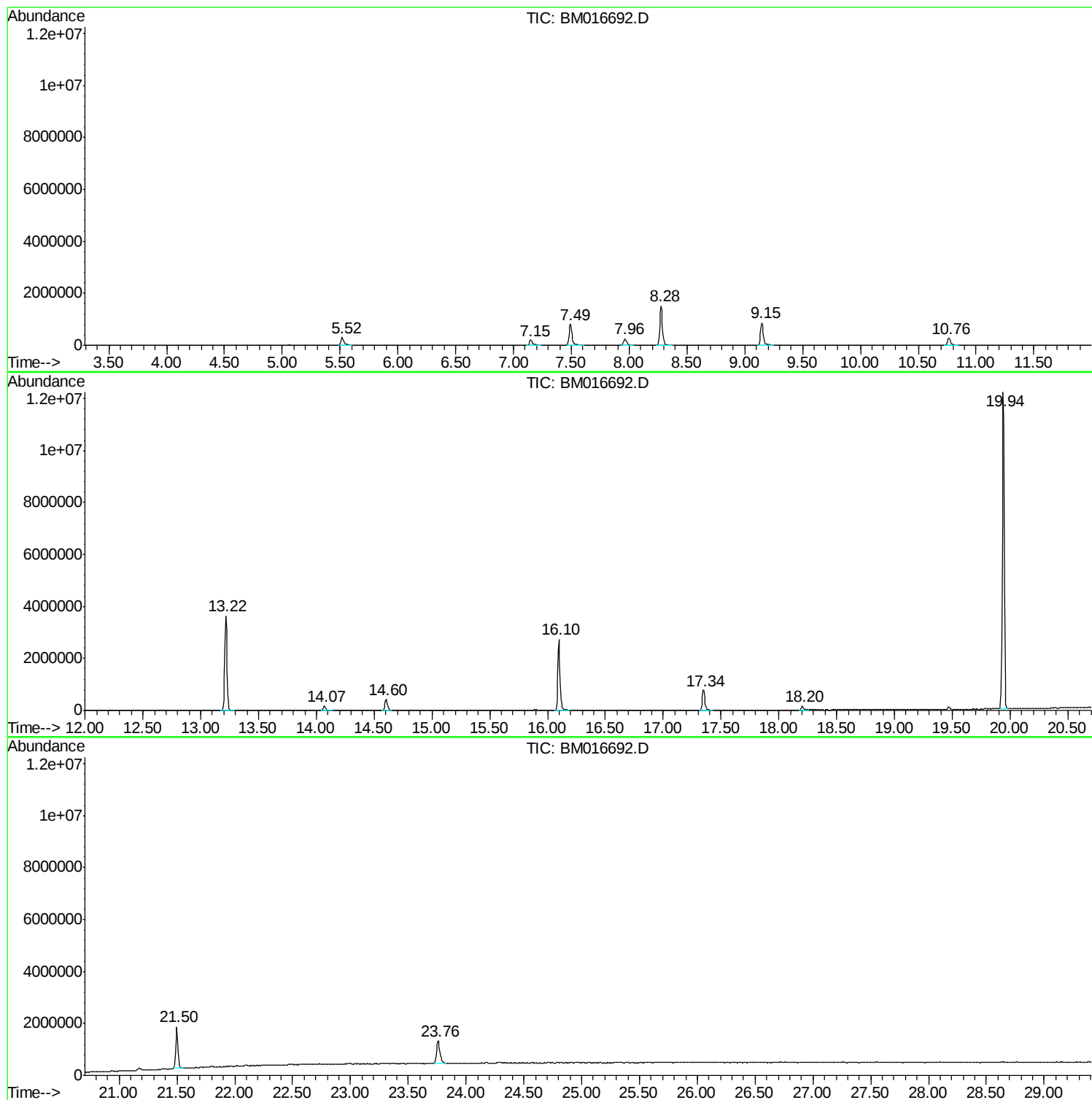
Sum of corrected areas: 36035145

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Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM082118.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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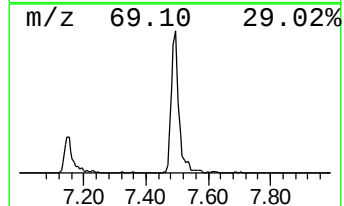
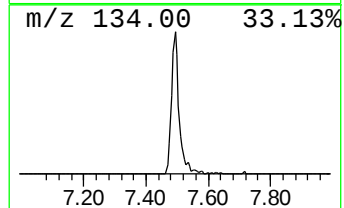
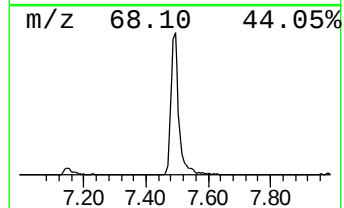
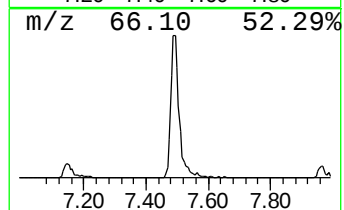
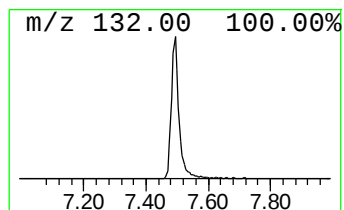
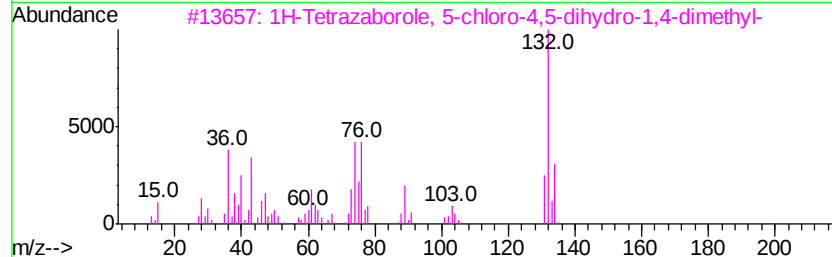
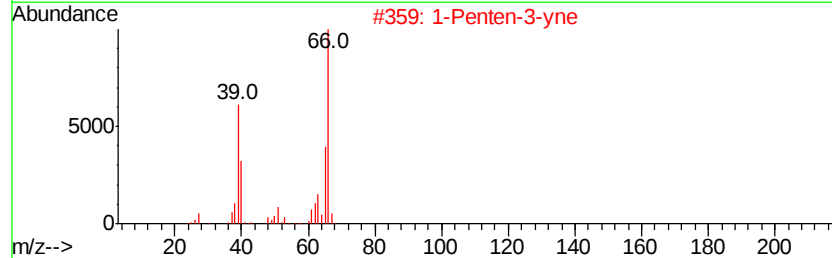
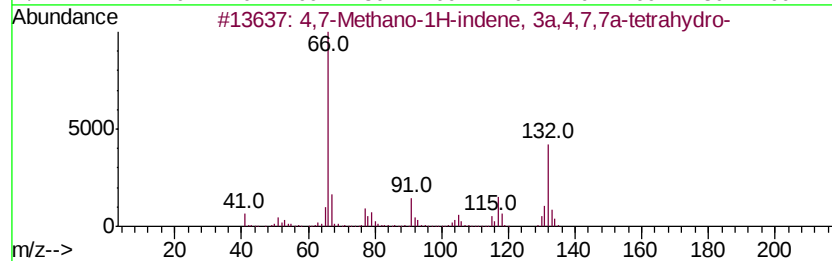
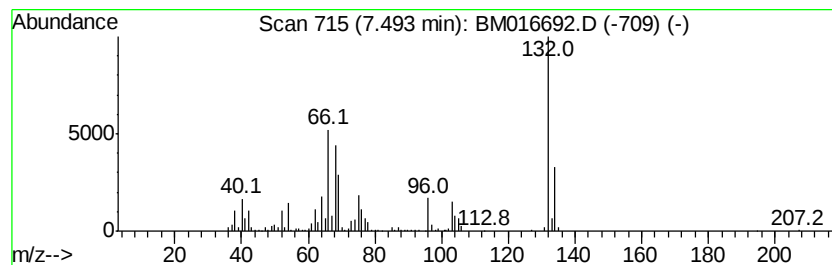
Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM082118.M
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 unknown7.49 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.49	61.42 ng	1436730	1,4-Dichlorobenzene-d4	7.96

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	4,7-Methano-1H-indene, 3a,4,7,7a...	132	C10H12	000077-73-6	38
2		1-Penten-3-yne	66	C5H6	000646-05-9	27
3		1H-Tetrazaborole, 5-chloro-4,5-d...	132	C2H6BClN4	021960-49-6	22
4		3-Penten-1-yne, (E)-	66	C5H6	002004-69-5	16
5		3-Penten-1-yne, (Z)-	66	C5H6	001574-40-9	16



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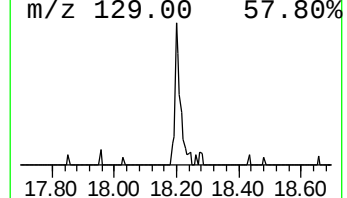
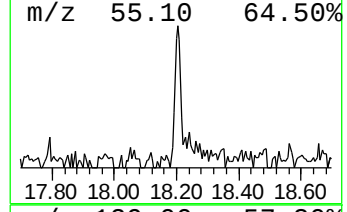
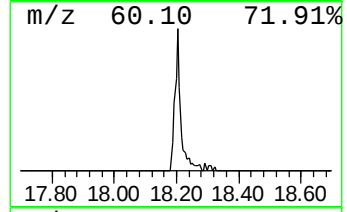
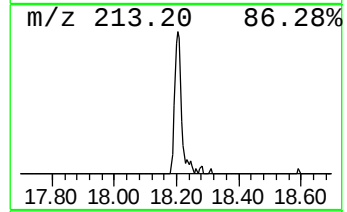
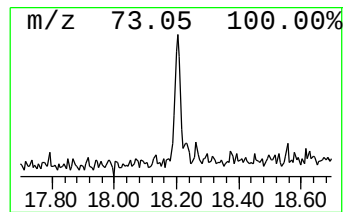
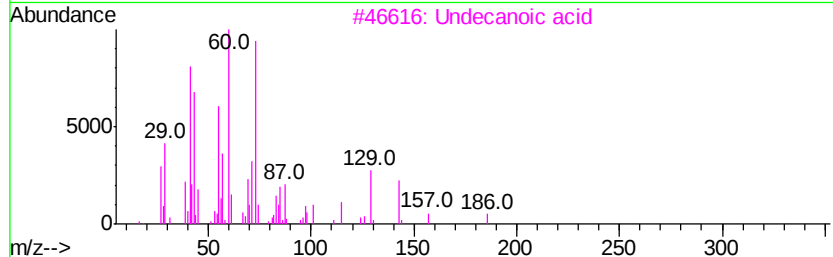
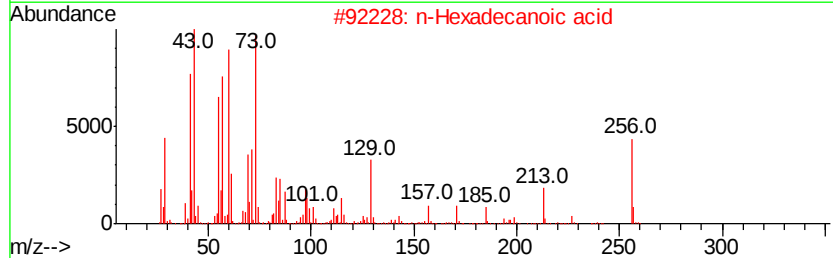
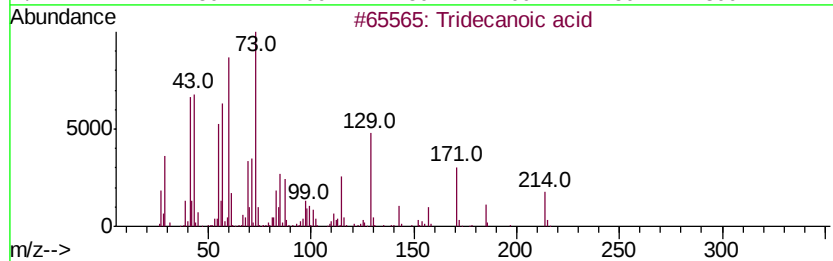
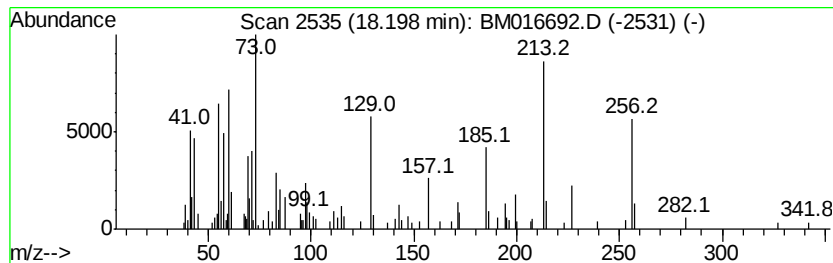
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Peak Number 3 unknown18.20 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.20	3.87 ng	226445	Phenanthrene-d10	17.34

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tridecanoic acid	214	C13H26O2	000638-53-9	46
2		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	45
3		Undecanoic acid	186	C11H22O2	000112-37-8	35
4		Lumiflavine	256	C13H12N4O2	001088-56-8	30
5		Tetradecanoic acid	228	C14H28O2	000544-63-8	27



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TIC Top Hit name	RT	EstConc	Units	Response	---Internal Standard---			
					#	RT	Resp	Conc
unknown7.49	7.49	61.4	ng	1436730	1	7.96	467841	20.0
unknown18.20	18.20	3.9	ng	226445	4	17.34	1170700	20.0