

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM090418\
 Data File : BM016658.D
 Acq On : 04 Sep 2018 22:19
 Operator : SJ/JU
 Sample : J4759-04
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 083018-SIDI-E-U-S

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	5.546	379	384	401	rBV	291230	593227	7.16%	2.055%
2	7.181	657	662	674	rBV	171591	370605	4.47%	1.284%
3	7.522	713	720	732	rBV2	811330	1428397	17.25%	4.948%
4	7.998	795	801	815	rBV2	285450	526943	6.36%	1.825%
5	8.310	847	854	869	rBV	1616424	2662375	32.15%	9.223%
6	9.181	996	1002	1025	rBV	861070	1587170	19.16%	5.498%
7	10.798	1272	1277	1293	rBV	292121	549837	6.64%	1.905%
8	13.251	1688	1694	1710	rBV	3577982	4893791	59.09%	16.953%
9	14.104	1833	1839	1846	rBV2	68132	113071	1.37%	0.392%
10	14.639	1922	1930	1938	rBV3	435090	722943	8.73%	2.504%
11	16.133	2178	2184	2199	rBV	1719822	2760152	33.33%	9.561%
12	17.380	2391	2396	2408	rBV	623207	960183	11.59%	3.326%
13	18.233	2536	2541	2551	rBV4	72501	105323	1.27%	0.365%
14	19.968	2831	2836	2845	rBV	7591394	8282223	100.00%	28.691%
15	21.527	3096	3101	3112	rBV	1069717	1518115	18.33%	5.259%
16	23.803	3482	3488	3496	rBV	850334	1793037	21.65%	6.211%

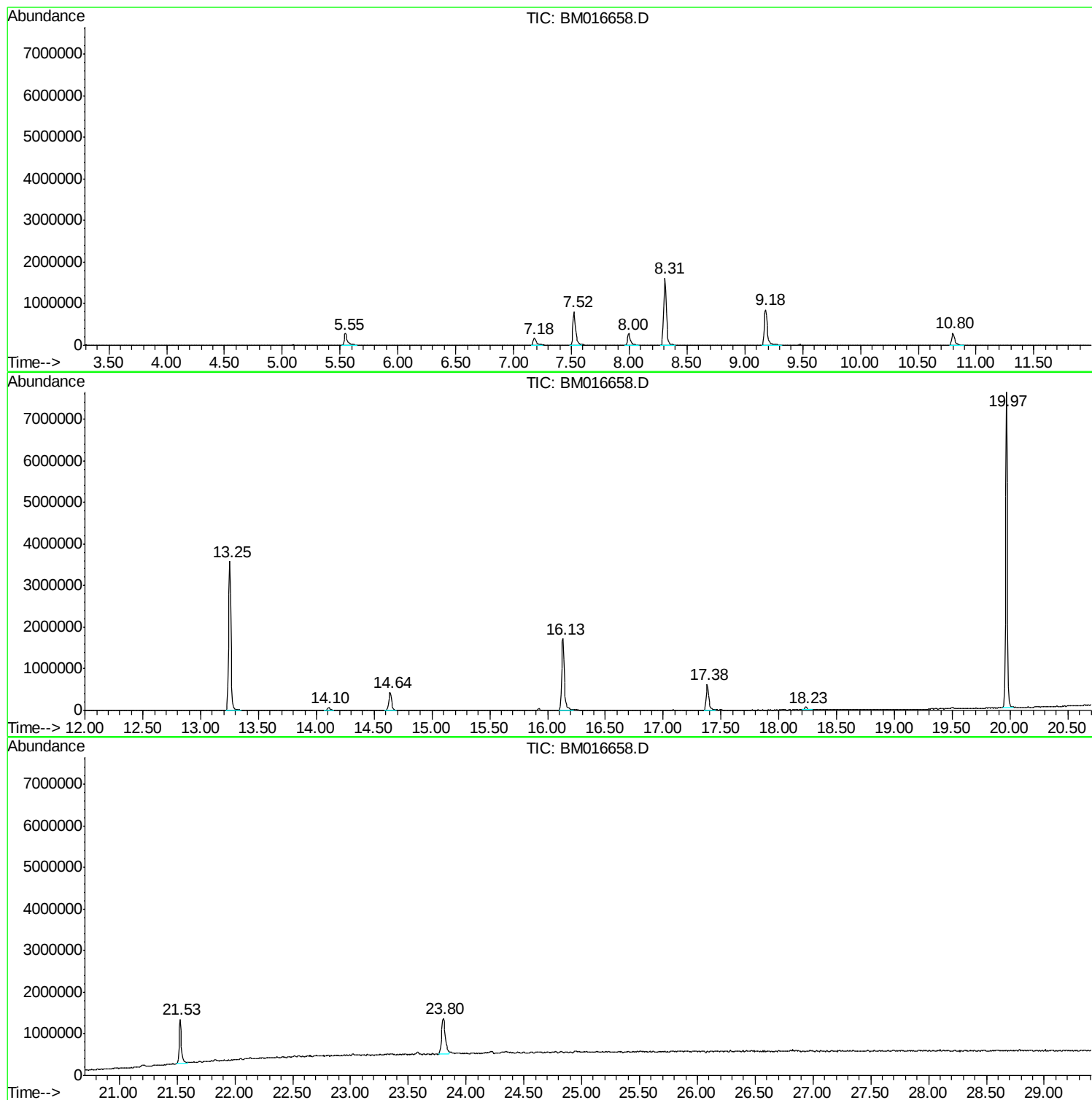
Sum of corrected areas: 28867392

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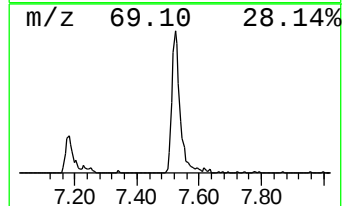
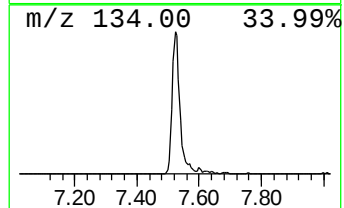
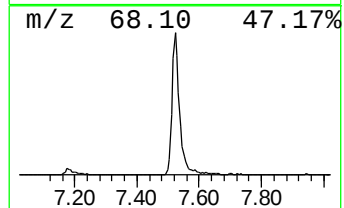
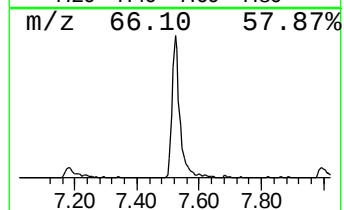
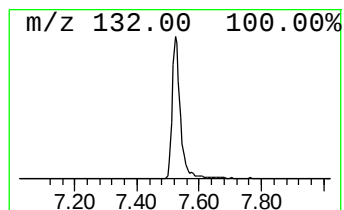
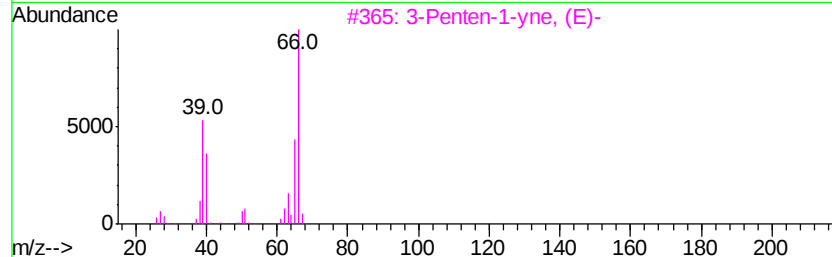
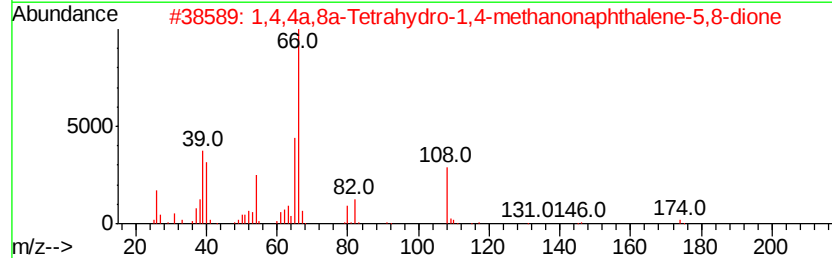
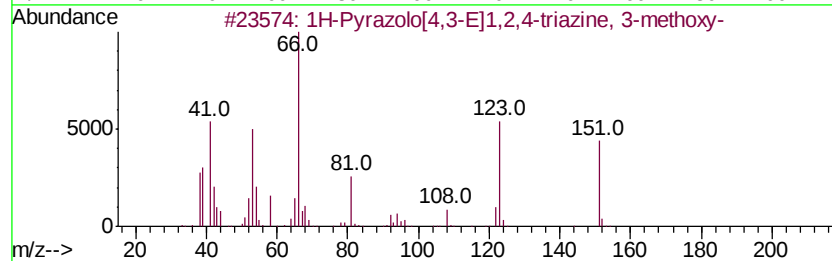
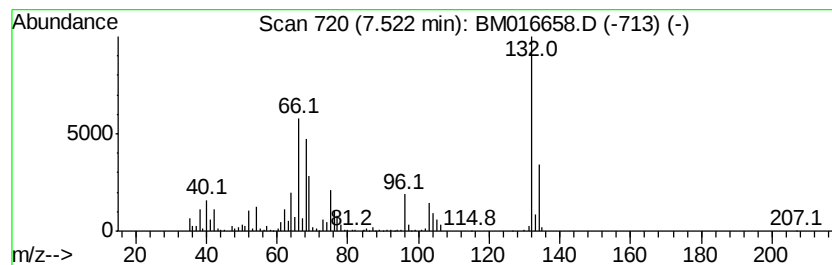
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 unknown7.52 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.52	54.21 ng	1428400	1,4-Dichlorobenzene-d4	7.99

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1H-Pyrazolo[4,3-E]1,2,4-triazine...	151	C5H5N5O	037526-53-7	16
2		1,4,4a,8a-Tetrahydro-1,4-methano...	174	C11H10O2	001200-89-1	16
3		3-Penten-1-yne, (E)-	66	C5H6	002004-69-5	16
4		1,3-Cyclopentadiene	66	C5H6	000542-92-7	16
5		1-Penten-3-yne	66	C5H6	000646-05-9	16



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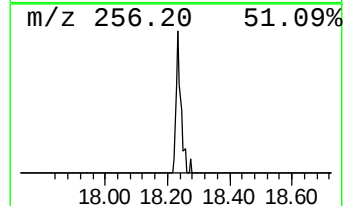
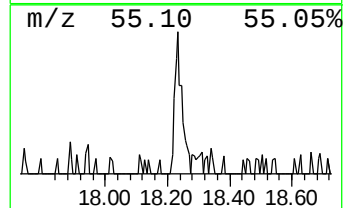
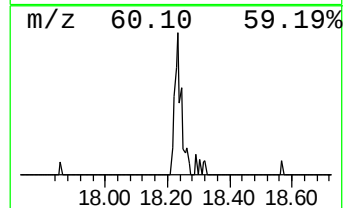
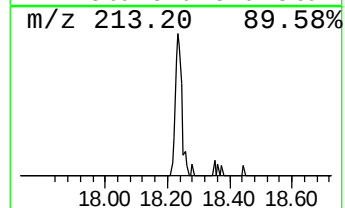
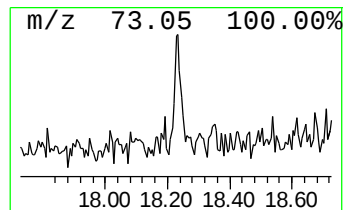
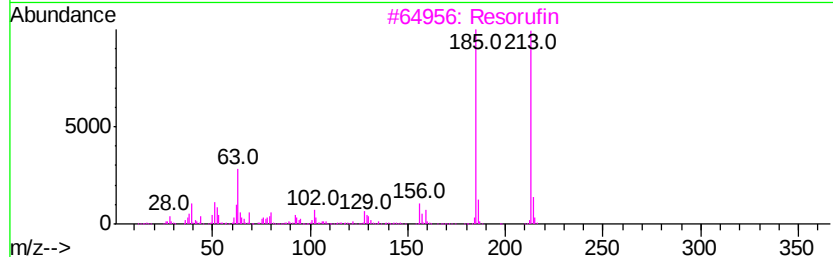
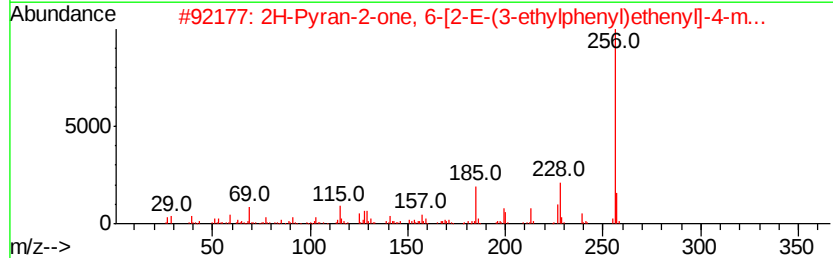
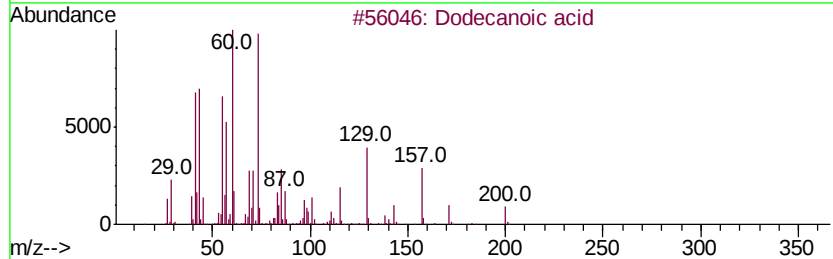
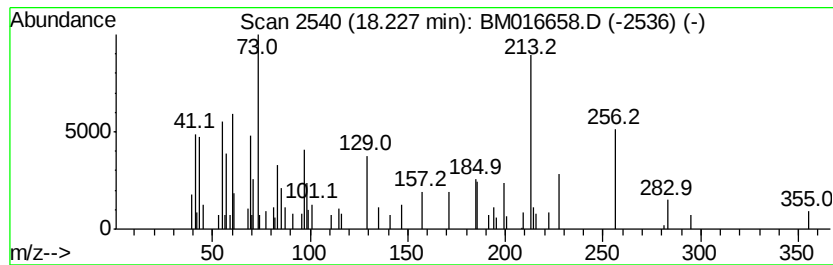
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Peak Number 3 Dodecanoic acid Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD		R.T.	
18.23	2.19 ng	105323	Phenanthrene-d10		17.38	
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Dodecanoic acid	200	C12H24O2	000143-07-7	64
2		2H-Pyran-2-one, 6-[2-E-(3-ethylp...	256	C16H16O3	1000159-82-7	46
3		Resorufin	213	C12H7NO3	000635-78-9	14
4		Tetradecanoic acid	228	C14H28O2	000544-63-8	14
5		Cyclopropane-1,2,3-tricarboxylic...	258	C12H18O6	013949-99-0	14



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TIC Top Hit name	RT	EstConc	Units	Response	---Internal Standard---			
					#	RT	Resp	Conc
unknown7.52	7.52	54.2	ng	1428400	1	7.99	526943	20.0
Dodecanoic acid	18.23	2.2	ng	105323	4	17.38	960183	20.0