

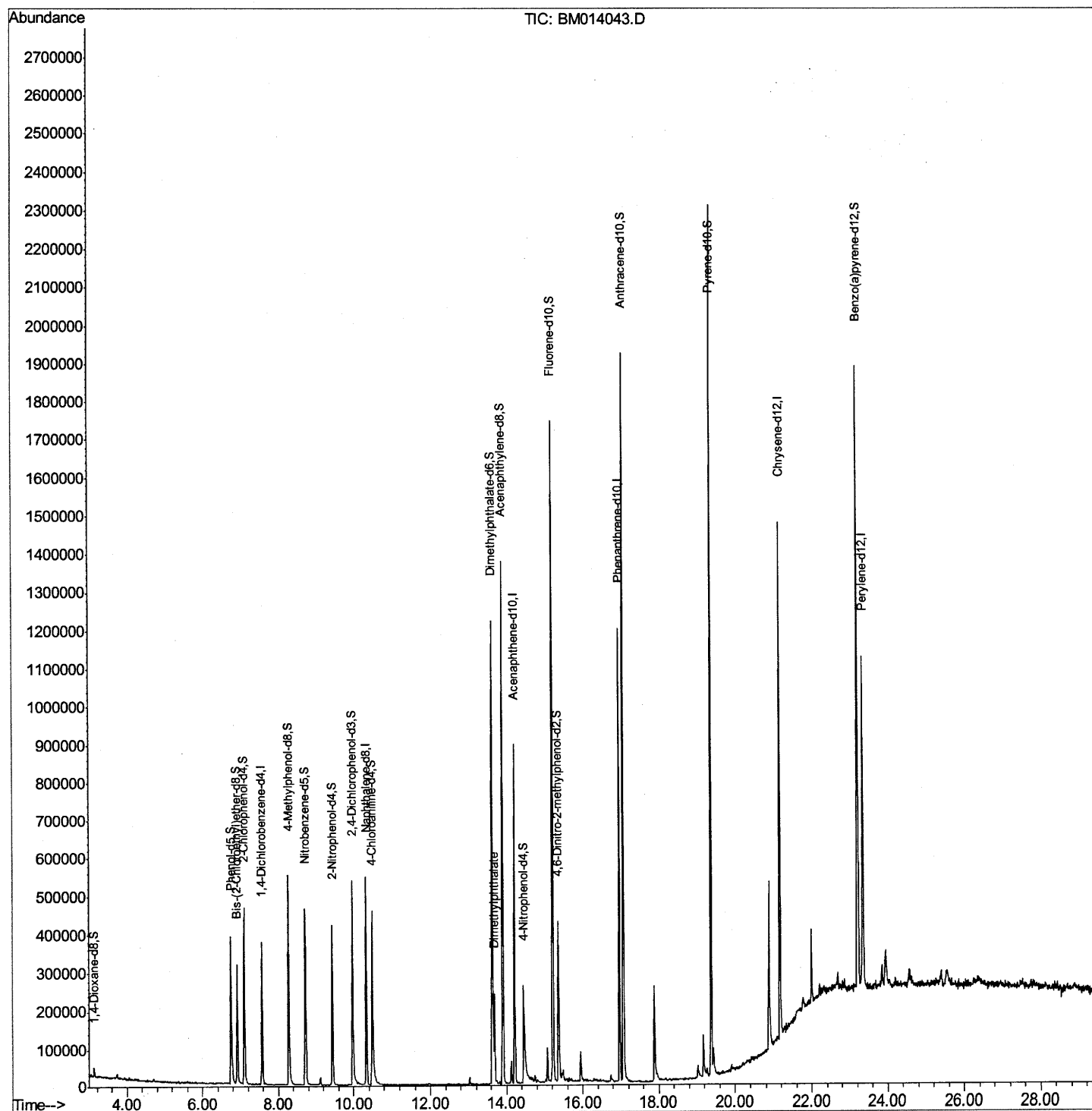
Data Path : Z:\HPCHEM1\BNA M\DATA\BM013118\
Data File : BM014043.D
Acq On : 31 Jan 2018 20:30
Operator : SJ/JU
Sample : J1314-18
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampled :
F6C55

Manual Integrations
APPROVED

Sohil
2/1/2018 6:04:43 PM

Quant Time: Feb 01 01:45:14 2018
Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM013118.M
Quant Title : SVOA CALIBRATION
QLast Update : Thu Feb 01 01:00:18 2018
Response via : Initial Calibration



Quantitation Report (Qedit)

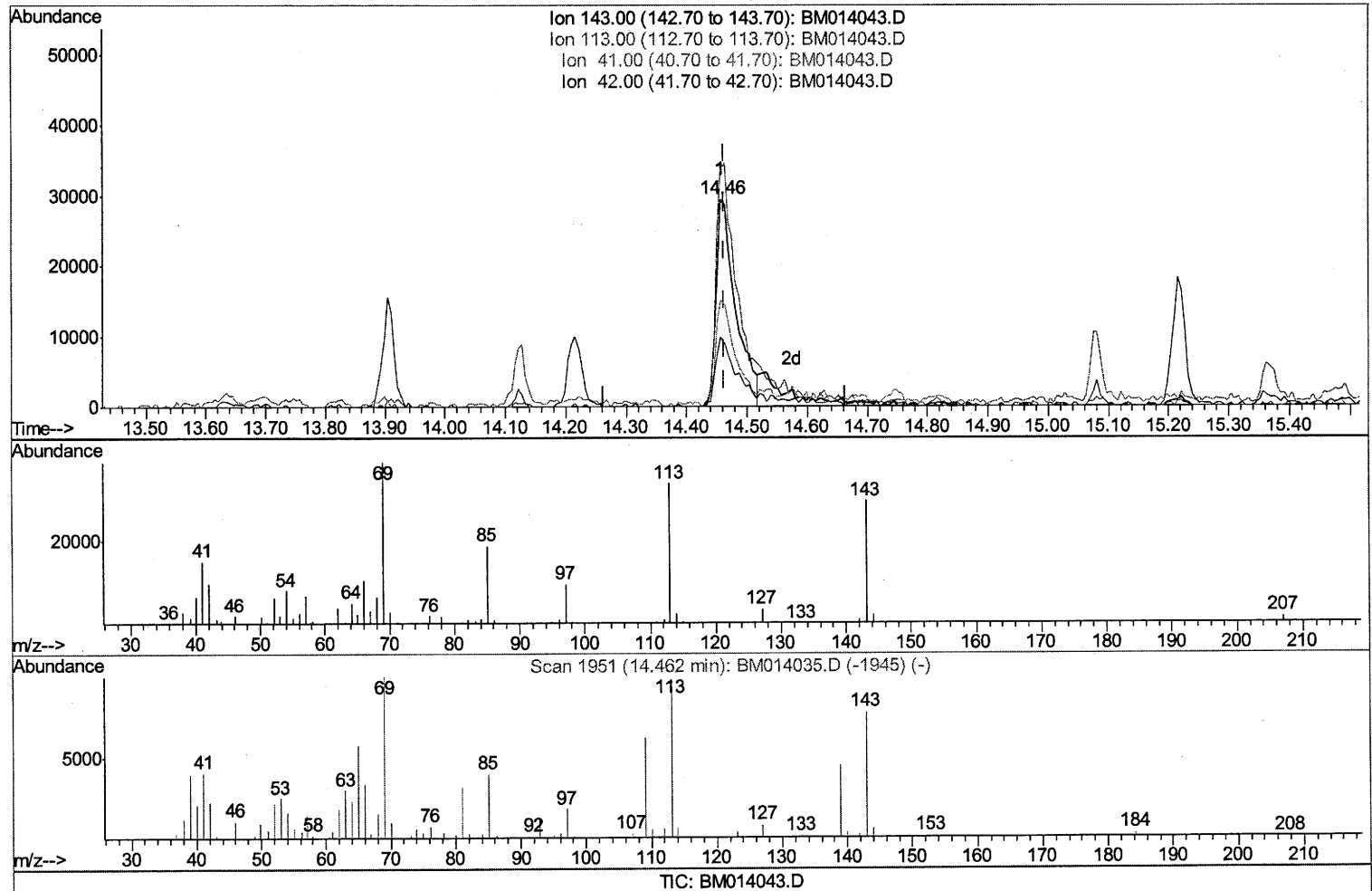
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Sohil
 2/1/2018 6:04:43 PM

Quant Time: Feb 01 01:04:05 2018
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM013118.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Feb 01 01:00:18 2018
 Response via : Initial Calibration



(51) 4-Nitrophenol-d4 (S)

14.457min (-0.006) 21.42ng/ul

response 65896

Ion	Exp%	Act%
143.00	100	100
113.00	114.60	114.37
41.00	34.80	51.13#
42.00	20.50	33.18#

Quantitation Report (Qedit)

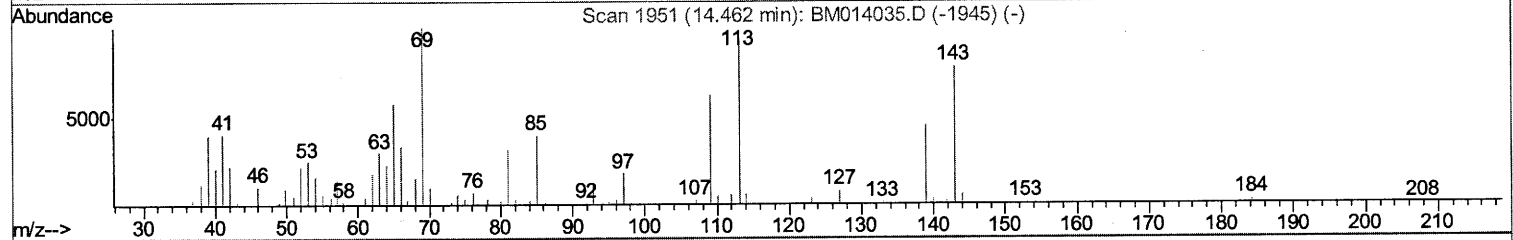
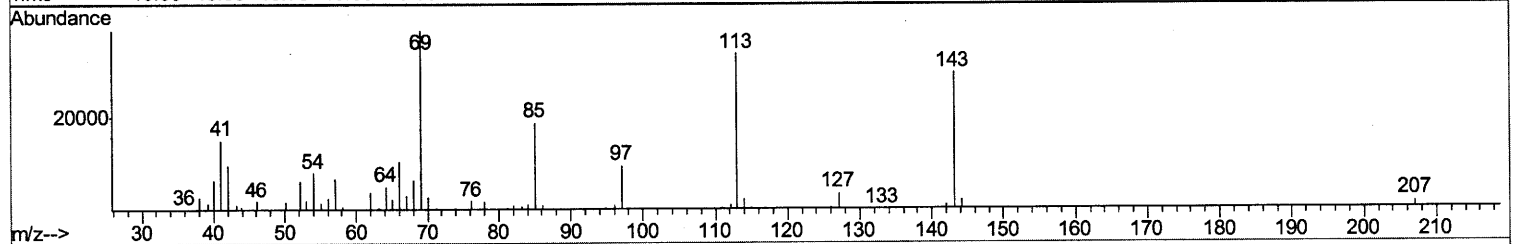
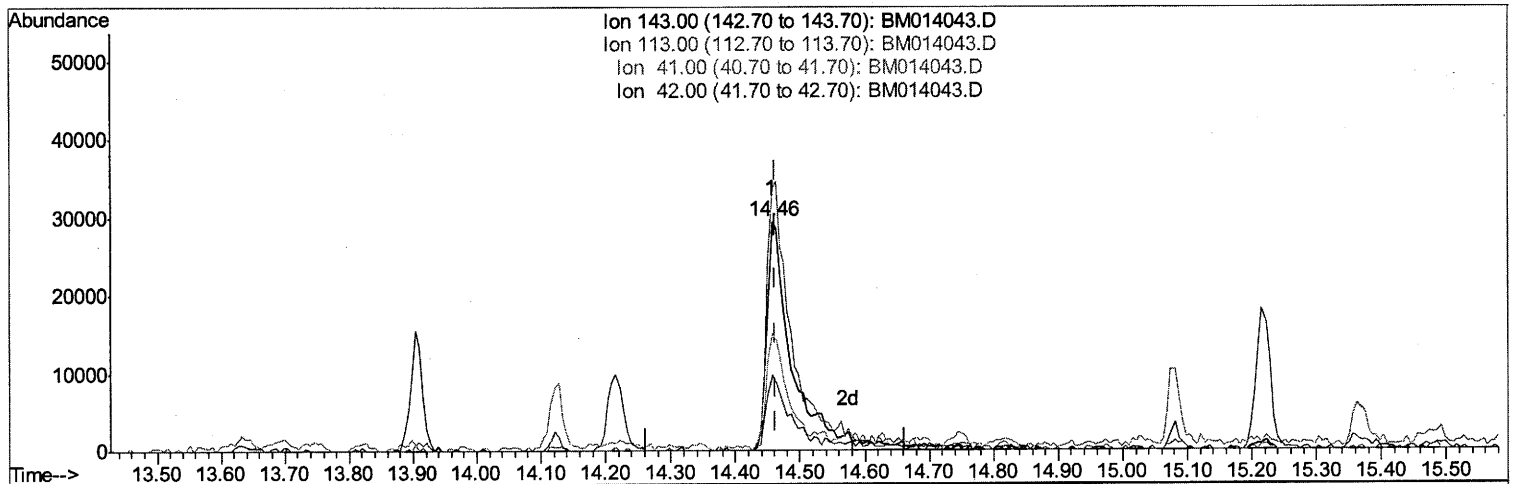
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 BNA_M
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 F6C55

Manual Integrations
 APPROVED

Sohil
 2/1/2018 6:04:43 PM

Quant Time: Feb 01 01:04:05 2018
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 Quant Title : SVOA CALIBRATION
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 Response via : Initial Calibration



TIC: BM014043.D

(51) 4-Nitrophenol-d4 (S)

14.457min (-0.006) 24.95ng/ul m > SJ 2/1/18

response 76767

Ion	Exp%	Act%
143.00	100	100
113.00	114.60	114.37
41.00	34.80	51.13#
42.00	20.50	33.18#

Data Path : Z:\HPCHEM1\BNA M\DATA\BM013118\
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 BNA_M
 Client Sampled :
 F6C55

Manual Integrations
 APPROVED

Sohil
 2/1/2018 6:04:43 PM

Quant Time: Feb 01 01:45:14 2018
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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.56	152	95295	20.00	ng/ul	0.00
18) Naphthalene-d8	10.33	136	392661	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.22	164	269033	20.00	ng/ul	0.00
61) Phenanthrene-d10	16.97	188	652560	20.00	ng/ul	0.00
75) Chrysene-d12	21.18	240	689193	20.00	ng/ul	0.00
83) Perylene-d12	23.36	264	654861	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.12	96	10820	5.21	ng/uL	0.00
5) Phenol-d5	6.75	99	202454	28.27	ng/ul	0.00
7) Bis-(2-Chloroethyl) ether-d	6.91	67	136522	31.60	ng/ul	0.00
9) 2-Chlorophenol-d4	7.10	132	166956	29.44	ng/ul	0.00
13) 4-Methylphenol-d8	8.27	113	188748	31.47	ng/ul	0.00
19) Nitrobenzene-d5	8.72	128	97210	34.30	ng/ul	0.00
22) 2-Nitrophenol-d4	9.43	143	105285	33.56	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	9.97	165	198945	30.15	ng/ul	0.00
29) 4-Chloroaniline-d4	10.49	131	232559	40.65	ng/ul	0.00
43) Dimethylphthalate-d6	13.64	166	763834	34.21	ng/ul	0.00
46) Acenaphthylene-d8	13.91	160	914797	33.77	ng/ul	0.00
51) 4-Nitrophenol-d4	14.46	143	76767m	24.95	ng/ul	0.00
57) Fluorene-d10	15.22	176	669127	33.49	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	15.36	200	98137	24.49	ng/ul	0.00
70) Anthracene-d10	17.07	188	1104661	34.87	ng/ul	0.00
76) Pyrene-d10	19.37	212	1218694	36.72	ng/ul	0.00
87) Benzo(a)pyrene-d12	23.22	264	1088167	33.86	ng/ul	0.00

S2 2/8/18

Target Compounds				Ovalue	
44) Dimethylphthalate	13.69	163	135862	6.312 ng/ul#	97

(#) = qualifier out of range (m) = manual integration (+) = signals summed