

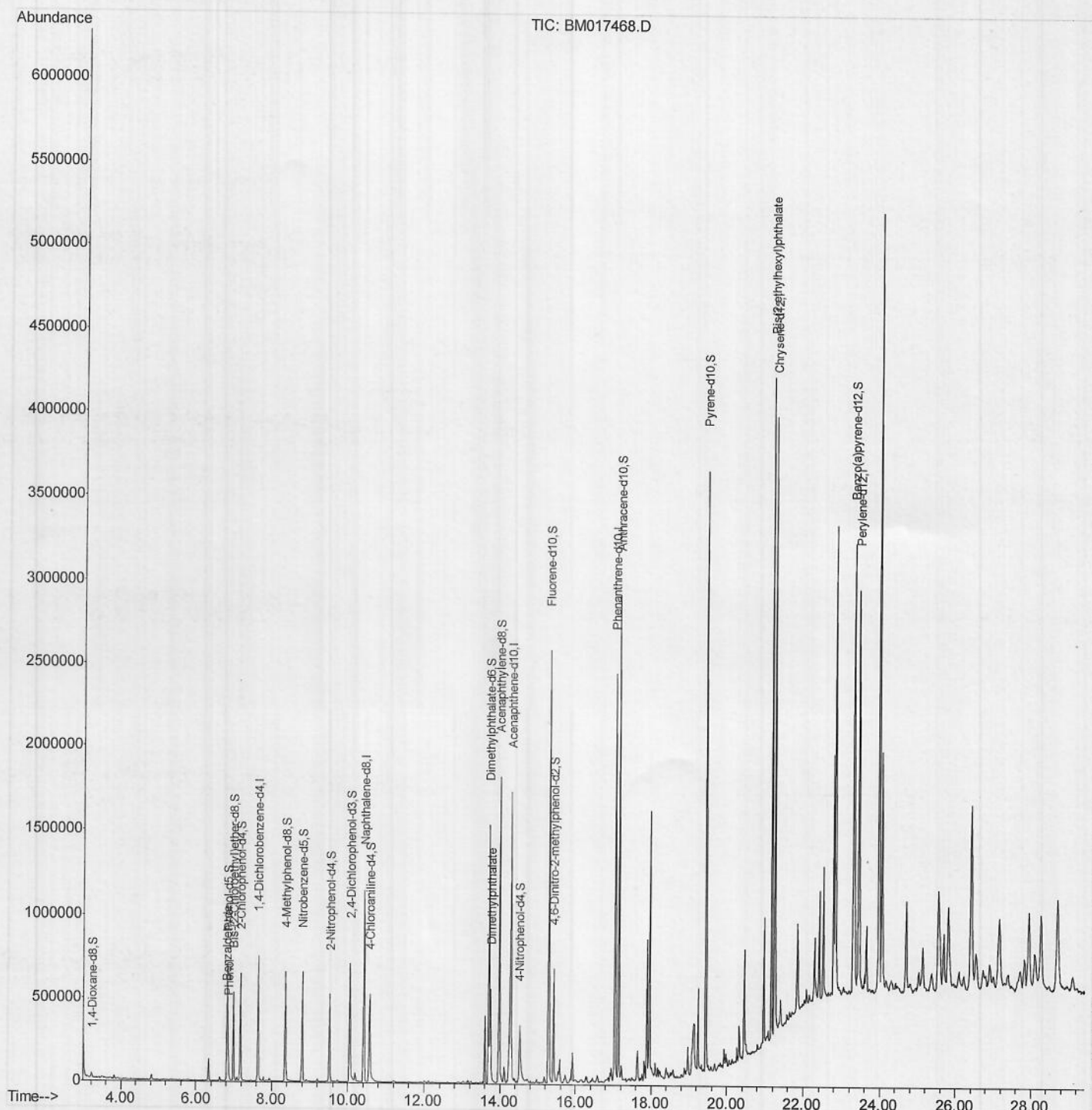
Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM110118\
 Quantitation Report (QT Reviewed)
 Data File : BM017468.D
 Acq On : 01 Nov 2018 16:46
 Operator : MJ/SJ
 Sample : J5704-17
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 A40H5

Manual Integrations
 APPROVED

Sohil
 11/2/2018 4:12:09 PM

Quant Time: Nov 02 02:55:06 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM102418MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Nov 02 02:41:22 2018
 Response via : Initial Calibration



Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM110118\
Quantitation Report (Qedit)

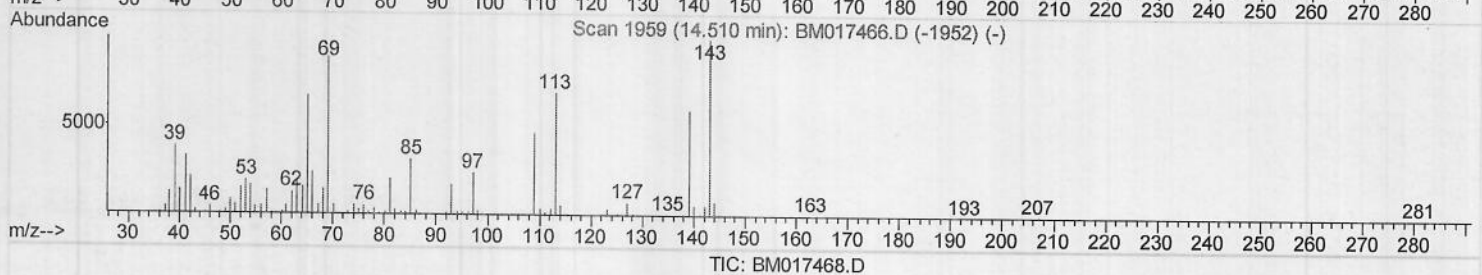
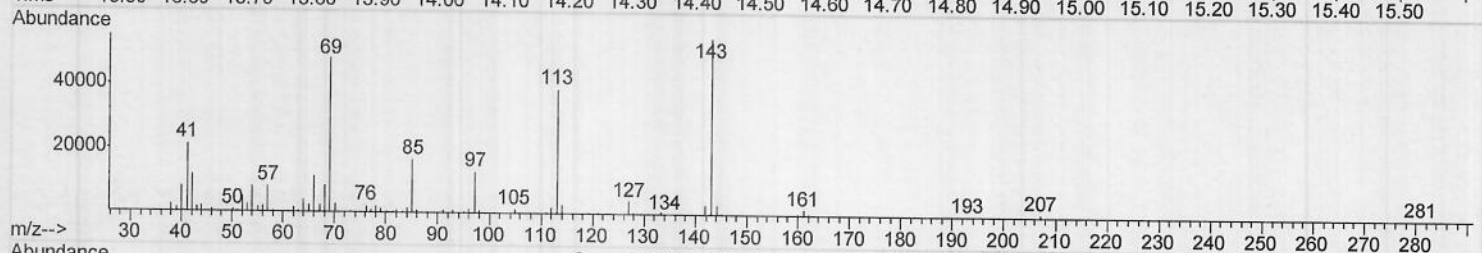
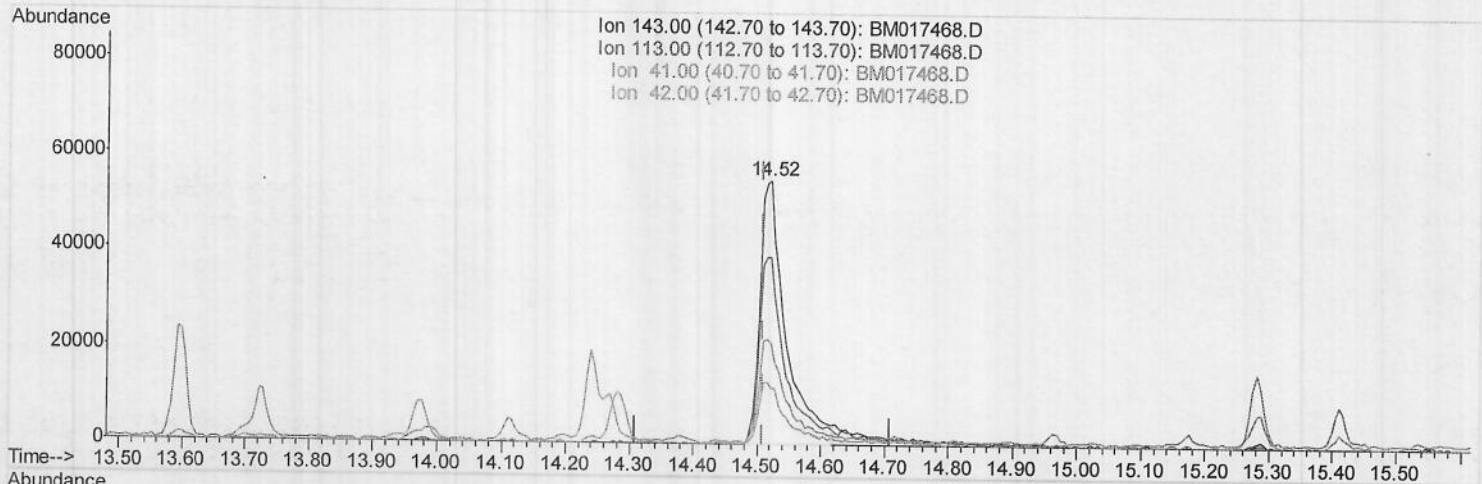
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Response via : Initial Calibration



(52) 4-Nitrophenol-d4 (S)

14.522min (+0.012) 16.07ng/ul

response 155089

Ion	Exp%	Act%
143.00	100	100
113.00	74.20	70.89
41.00	39.30	38.75
42.00	25.10	21.36

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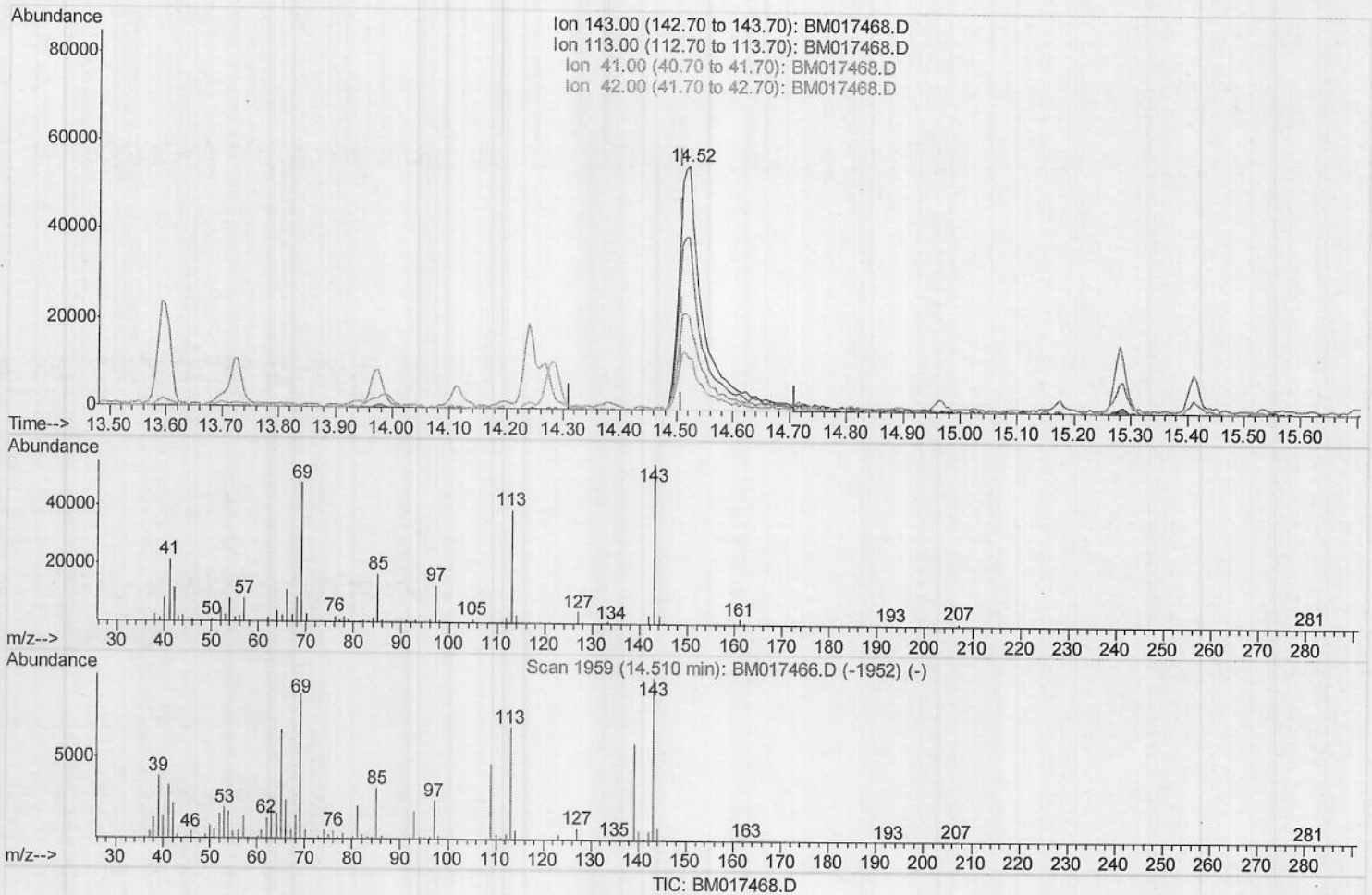
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Response via : Initial Calibration



(52) 4-Nitrophenol-d4 (S)

14.522min (+0.012) 17.34ng/ul m) SJ 11/06/18

response 167378

Ion	Exp%	Act%
143.00	100	100
113.00	74.20	70.89
41.00	39.30	38.75
42.00	25.10	21.36

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Manual Integrations
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Quant Title : SVOA CALIBRATION

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Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.64	152	201864	20.00	ng/ul	0.00
18) Naphthalene-d8	10.41	136	961435	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.28	164	602349	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.03	188	1463008	20.00	ng/ul	0.00
78) Chrysene-d12	21.23	240	1763733	20.00	ng/ul	0.00
86) Perylene-d12	23.44	264	1719436	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.22	96	16310	3.72	ng/uL	0.00
5) Phenol-d5	6.82	99	374727	20.32	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	6.99	67	225421	20.96	ng/ul	0.00
9) 2-Chlorophenol-d4	7.18	132	303178	21.21	ng/ul	0.00
13) 4-Methylphenol-d8	8.35	113	263241	18.06	ng/ul	0.00
19) Nitrobenzene-d5	8.79	128	160729	21.39	ng/ul	0.00
22) 2-Nitrophenol-d4	9.51	143	180949	21.47	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.04	165	311050	19.87	ng/ul	0.00
29) 4-Chloroaniline-d4	10.56	131	354497	27.74	ng/ul	0.00
44) Dimethylphthalate-d6	13.70	166	1110524	21.50	ng/ul	0.00
47) Acenaphthylene-d8	13.97	160	1382995	22.05	ng/ul	0.00
52) 4-Nitrophenol-d4	14.52	143	167378m	17.34	ng/ul	0.01
58) Fluorene-d10	15.28	176	1003924	22.38	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.41	200	153952	15.10	ng/ul	0.00
71) Anthracene-d10	17.13	188	1582782	21.86	ng/ul	0.00
79) Pyrene-d10	19.43	212	1891146	23.14	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.30	264	2046753	22.54	ng/ul	0.00

Target Compounds

					Qvalue
4) Benzaldehyde	6.80	77	6648	1.892	ng/ul 89
6) Phenol	6.85	94	99244	5.420	ng/ul 98
45) Dimethylphthalate	13.75	163	364375	7.281	ng/ul 100
84) Bis(2-ethylhexyl)phthalate	21.16	149	1360341	22.768	ng/ul 99

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

5/11/06/18