

Quantitation Report (QT Reviewed)

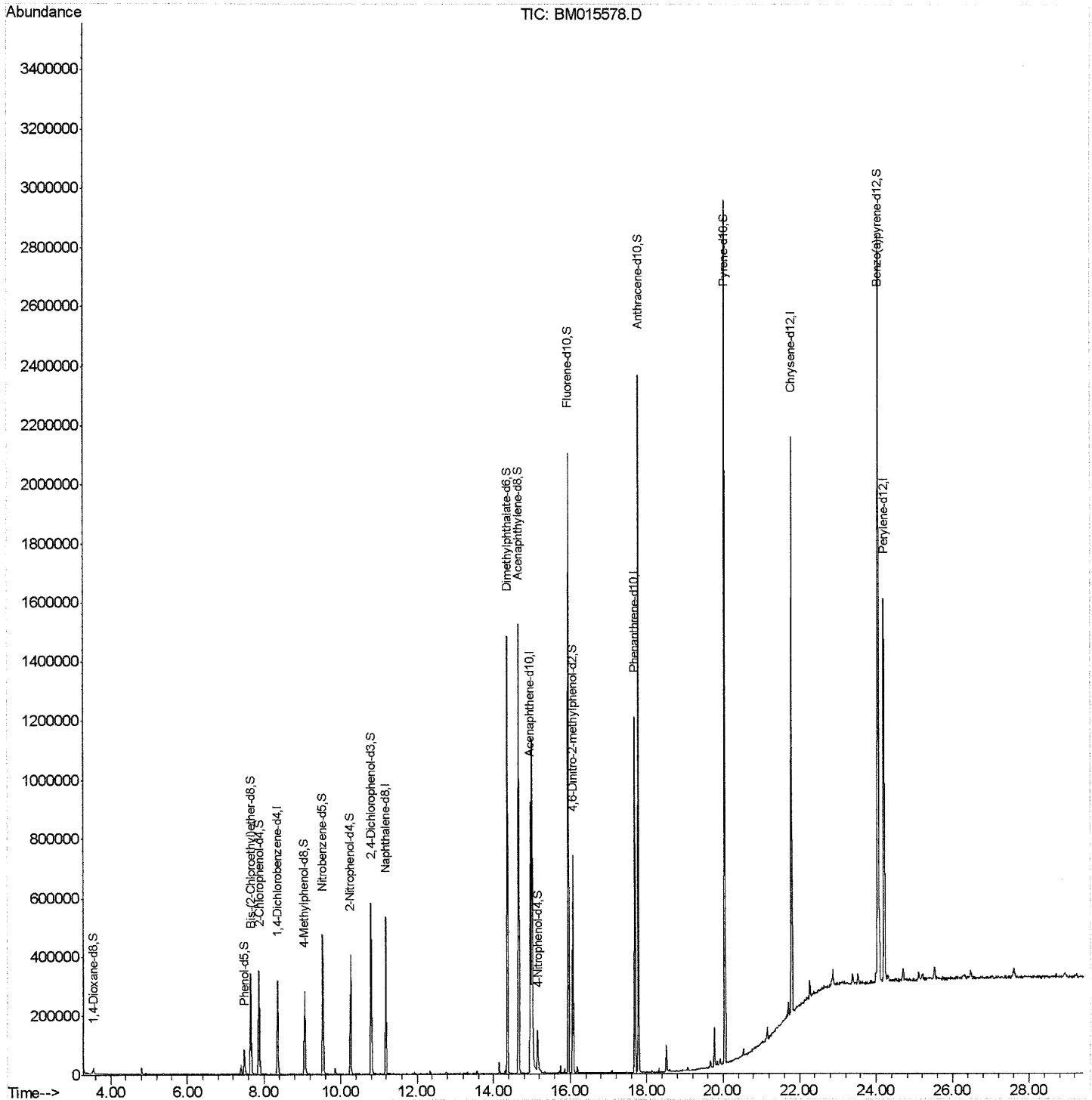
Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM060818\
 Data File : BM015578.D
 Acq On : 09 Jun 2018 00:04
 Operator : SJ/JU
 Sample : J3345-03
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampled :
 DAXF2

Manual Integrations
 APPROVED

Sohil
 6/11/2018 3:59:17 PM

Quant Time: Jun 09 02:16:15 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM060818MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Jun 08 23:05:23 2018
 Response via : Initial Calibration



Quantitation Report (Qedit)

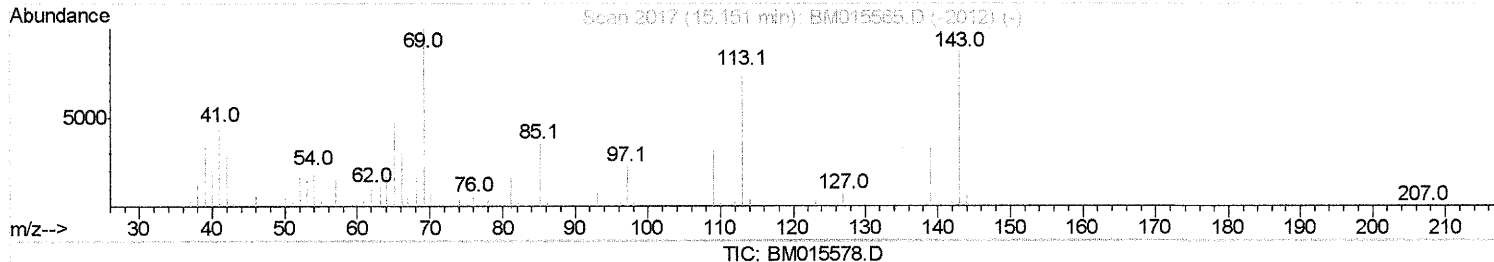
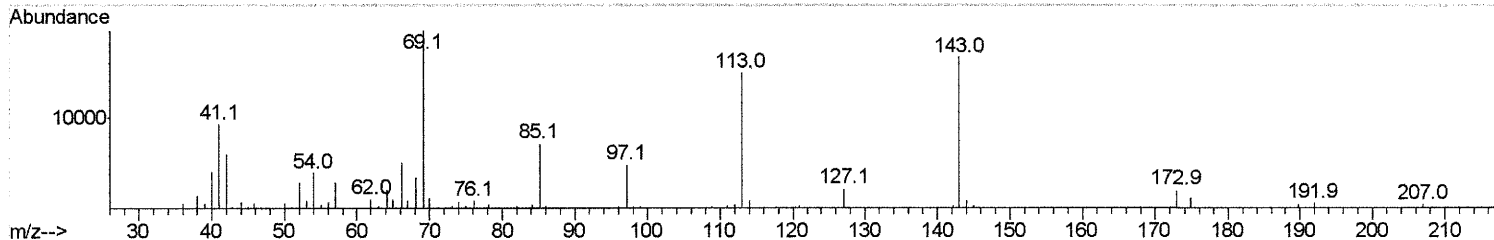
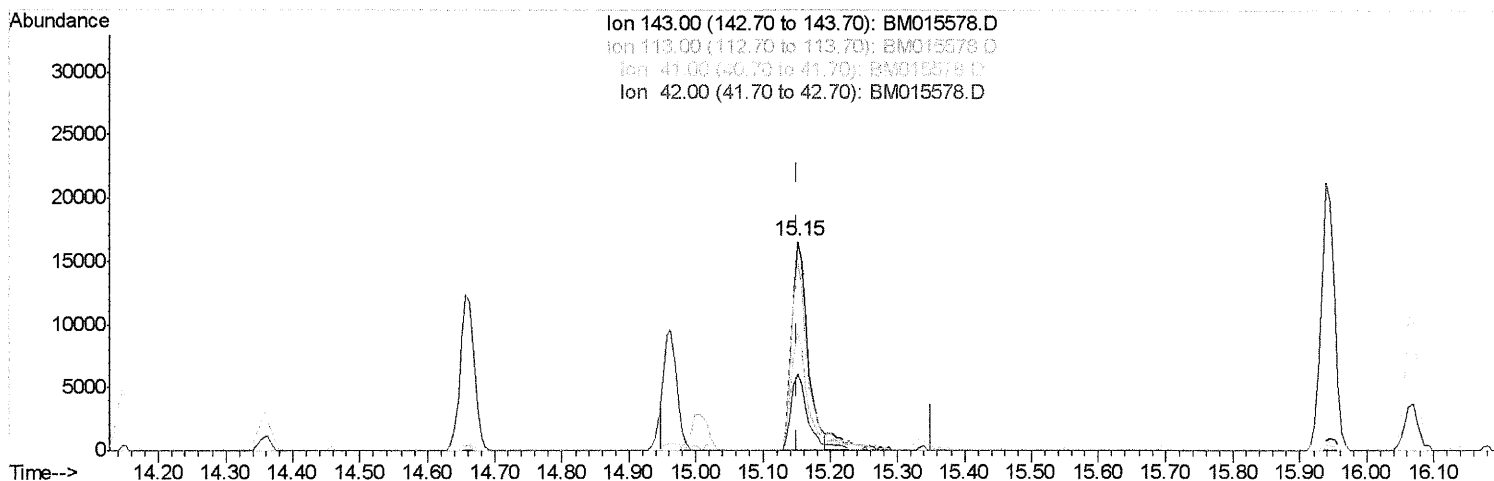
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TIC: BM015578.D

(51) 4-Nitrophenol-d4 (S)

15.151min (+0.000) 6.22ng/ul

response 27247

Ion	Exp%	Act%
143.00	100	100
113.00	75.10	89.95
41.00	28.80	56.37#
42.00	19.10	37.04#

Quantitation Report (Qedit)

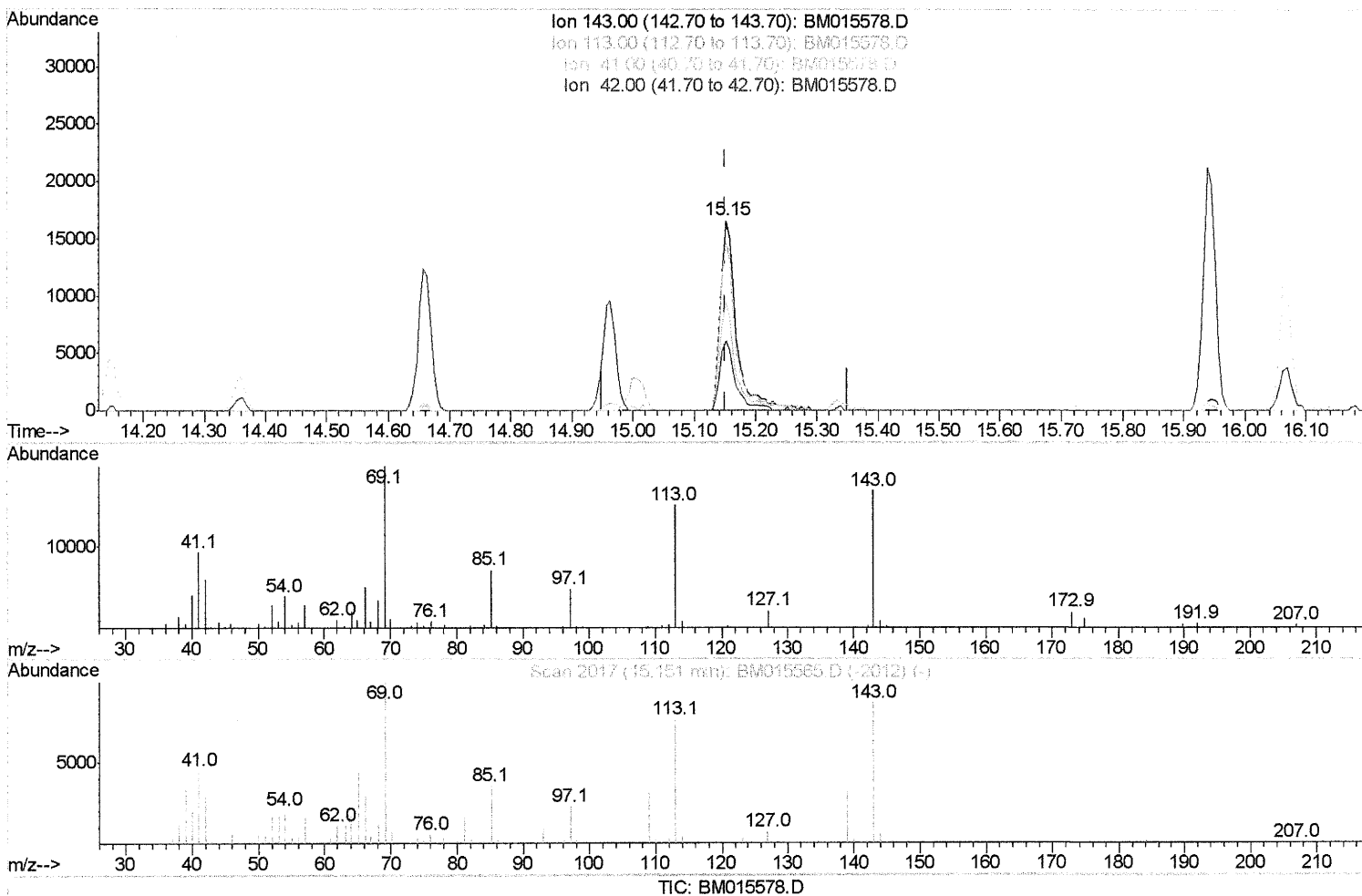
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 Operator : SJ/JU
 Sample : J3345-03
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampled :
 DAXF2

Manual Integrations
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(51) 4-Nitrophenol-d4 (S)

15.151min (+0.000) 6.92ng/ul m

SJ 6/11/18

response 30319

Ion	Exp%	Act%
143.00	100	100
113.00	75.10	89.95
41.00	28.80	56.37#
42.00	19.10	37.04#

Quantitation Report (QT Reviewed)

Instrument :
BNA_M
ClientSampleId :
DAXF2

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM060818\
Data File : BM015578.D
Acq On : 09 Jun 2018 00:04
Operator : SJ/JU
Sample : J3345-03
Misc :
ALS Vial : 22 Sample Multiplier: 1

Manual Integrations
APPROVED

Sohil
6/11/2018 3:59:17 PM

Quant Time: Jun 09 02:16:15 2018
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Quant Title : SVOA CALIBRATION
QLast Update : Fri Jun 08 23:05:23 2018
Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.35	152	81293	20.00	ng/ul	0.00
18) Naphthalene-d8	11.17	136	410618	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.96	164	274285	20.00	ng/ul	0.00
61) Phenanthrene-d10	17.69	188	669733	20.00	ng/ul	0.00
77) Chrysene-d12	21.79	240	836463	20.00	ng/ul	0.00
85) Perylene-d12	24.20	264	943748	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.55	96	10783	6.45	ng/uL	0.00
5) Phenol-d5	7.49	99	47372	6.42	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.66	67	158518	35.55	ng/ul	0.00
9) 2-Chlorophenol-d4	7.87	132	154471	27.37	ng/ul	0.00
13) 4-Methylphenol-d8	9.05	113	103869	16.34	ng/ul	0.00
19) Nitrobenzene-d5	9.53	128	101470	36.75	ng/ul	0.00
22) 2-Nitrophenol-d4	10.26	143	112243	37.80	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.79	165	196282	31.92	ng/ul	0.00
29) 4-Chloroaniline-d4	0.00	131	0d	0.00	ng/ul	0.00
43) Dimethylphthalate-d6	14.36	166	887360	39.08	ng/ul	0.00
46) Acenaphthylene-d8	14.66	160	1003910	40.99	ng/ul	0.00
51) 4-Nitrophenol-d4	15.15	143	30319m	6.92	ng/ul	0.00
57) Fluorene-d10	15.94	176	762637	39.69	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	16.07	200	129776	32.17	ng/ul	0.00
70) Anthracene-d10	17.78	188	1201932	39.37	ng/ul	0.00
78) Pyrene-d10	20.03	212	1376166	38.25	ng/ul	0.00
89) Benzo(a)pyrene-d12	24.04	264	1881396	39.77	ng/ul	0.00

316/11/18

Target Compounds Ovalue

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