

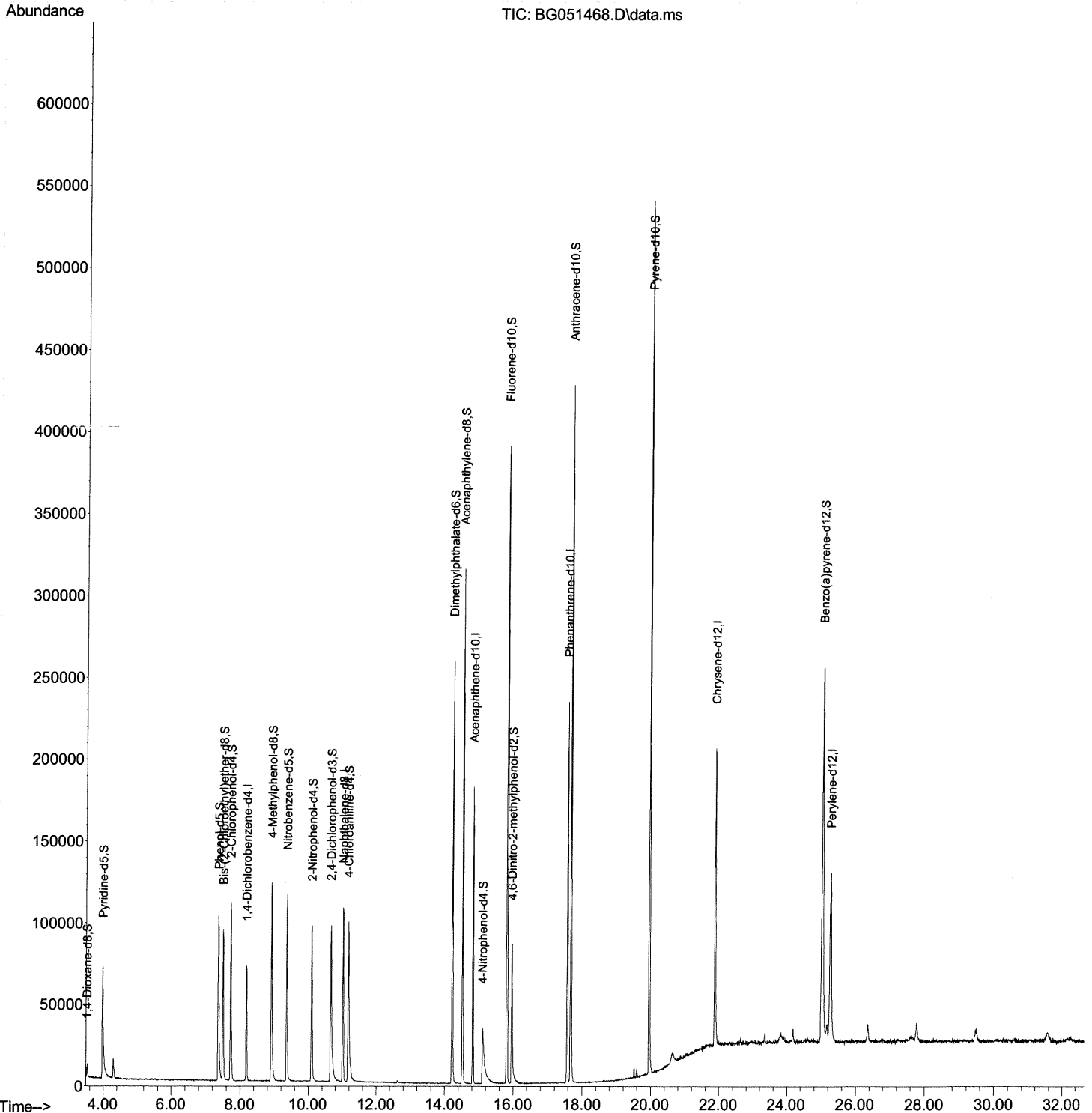
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120921\
Data File : BG051468.D
Acq On : 10 Dec 2021 23:46
Operator : CG/JU
Sample : PB141311BL
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
SBLK311

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 12/13/2021
Supervised By :Yogesh Patel 12/15/2021

Quant Time: Dec 11 04:26:18 2021
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG120821.M
Quant Title : SVOA CALIBRATION
QLast Update : Thu Dec 09 03:21:41 2021
Response via : Initial Calibration



Quantitation Report (Qedit)

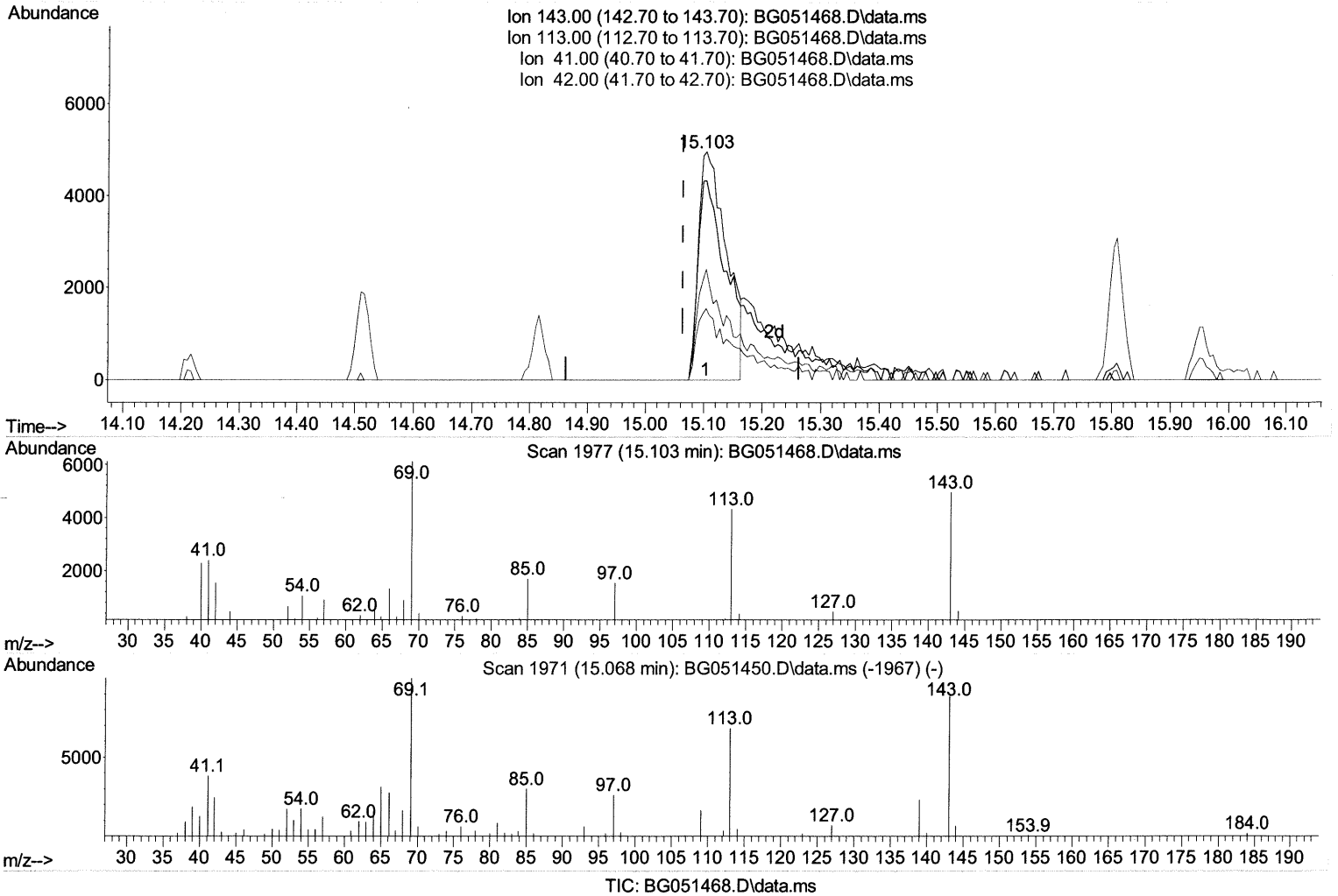
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(54) 4-Nitrophenol-d4 (S)

15.103min (+ 0.040) 22.35 ng/ul

response 16542

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	80.30	87.25
41.00	44.40	48.54
42.00	29.70	31.24

Quantitation Report (Qedit)

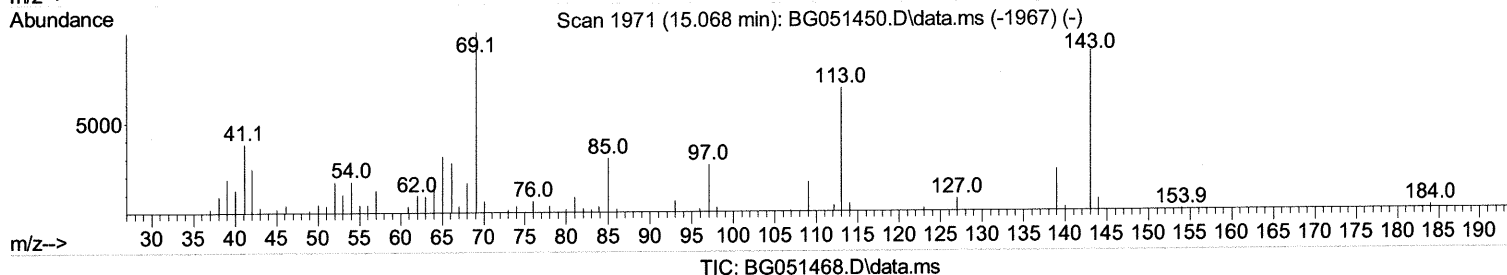
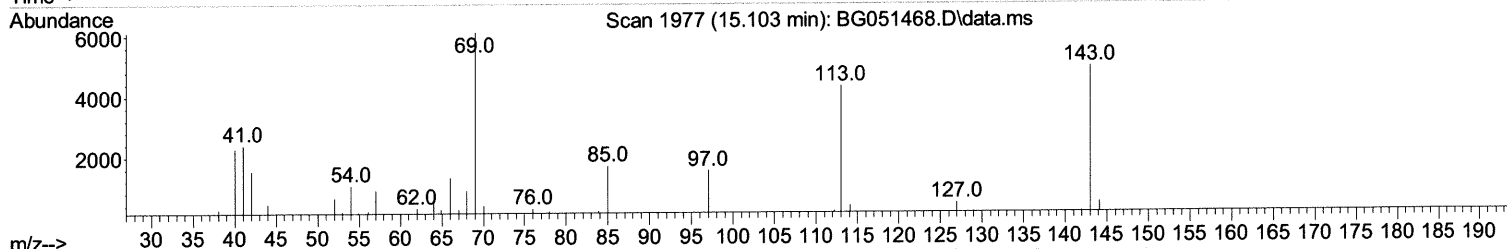
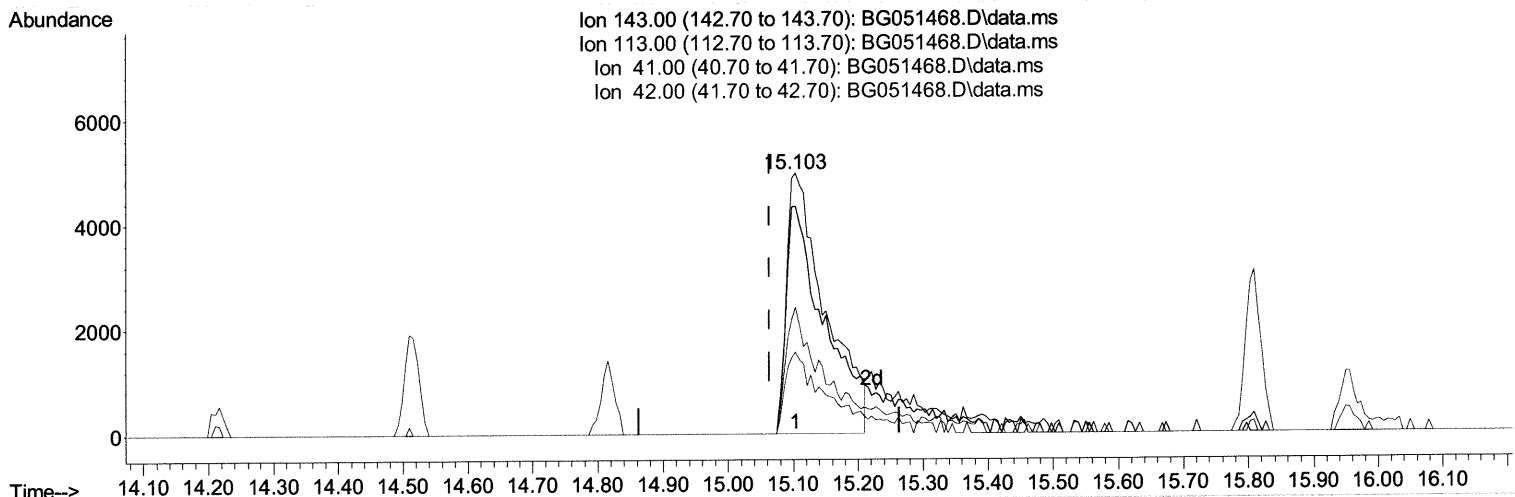
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TIC: BG051468.D\data.ms

(54) 4-Nitrophenol-d4 (S)

15.103min (+ 0.040) 27.73 ng/ul m 12/16/21 JU

response 20531

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	80.30	87.25
41.00	44.40	48.54
42.00	29.70	31.24

Quantitation Report (QT Reviewed)

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 Data File : BG051468.D
 Acq On : 10 Dec 2021 23:46
 Operator : CG/JU
 Sample : PB141311BL
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SBLK311

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 12/13/2021
 Supervised By :Yogesh Patel 12/15/2021

Quant Time: Dec 11 04:26:18 2021
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.182	152	21407	20.000	ng/ul	0.00
20) Naphthalene-d8	11.008	136	93459	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.815	164	63542	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.565	188	148021	20.000	ng/ul	0.00
79) Chrysene-d12	21.872	240	136233	20.000	ng/ul	0.00
88) Perylene-d12	25.268	264	123033	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.529	96	4308	6.609	ng/uL	0.00
4) Pyridine-d5	3.975	84	53214	28.428	ng/ul	0.01
7) Phenol-d5	7.365	99	74739	34.295	ng/ul	0.01
9) Bis-(2-Chloroethyl)eth...	7.500	67	49513	35.429	ng/ul	0.00
11) 2-Chlorophenol-d4	7.724	132	55176	35.587	ng/ul	0.00
15) 4-Methylphenol-d8	8.916	113	57263	33.448	ng/ul	0.00
21) Nitrobenzene-d5	9.369	128	29922	36.908	ng/ul	0.00
24) 2-Nitrophenol-d4	10.097	143	33472	36.486	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.656	165	51514	34.516	ng/ul	0.00
31) 4-Chloroaniline-d4	11.167	131	75045	34.378	ng/ul	0.00
46) Dimethylphthalate-d6	14.216	166	191144	38.876	ng/ul	0.00
49) Acenaphthylene-d8	14.516	160	236049	37.906	ng/ul	0.00
54) 4-Nitrophenol-d4	15.103	143	20531m	27.735	ng/ul	0.04
60) Fluorene-d10	15.808	176	164945	37.686	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.955	200	25170	28.615	ng/ul	0.00
73) Anthracene-d10	17.665	188	274698	39.663	ng/ul	0.00
81) Pyrene-d10	19.951	212	328971	40.175	ng/ul	0.00
92) Benzo(a)pyrene-d12	25.027	264	270742	42.664	ng/ul	0.00

Target Compounds Qvalue

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