

(QT Reviewed)

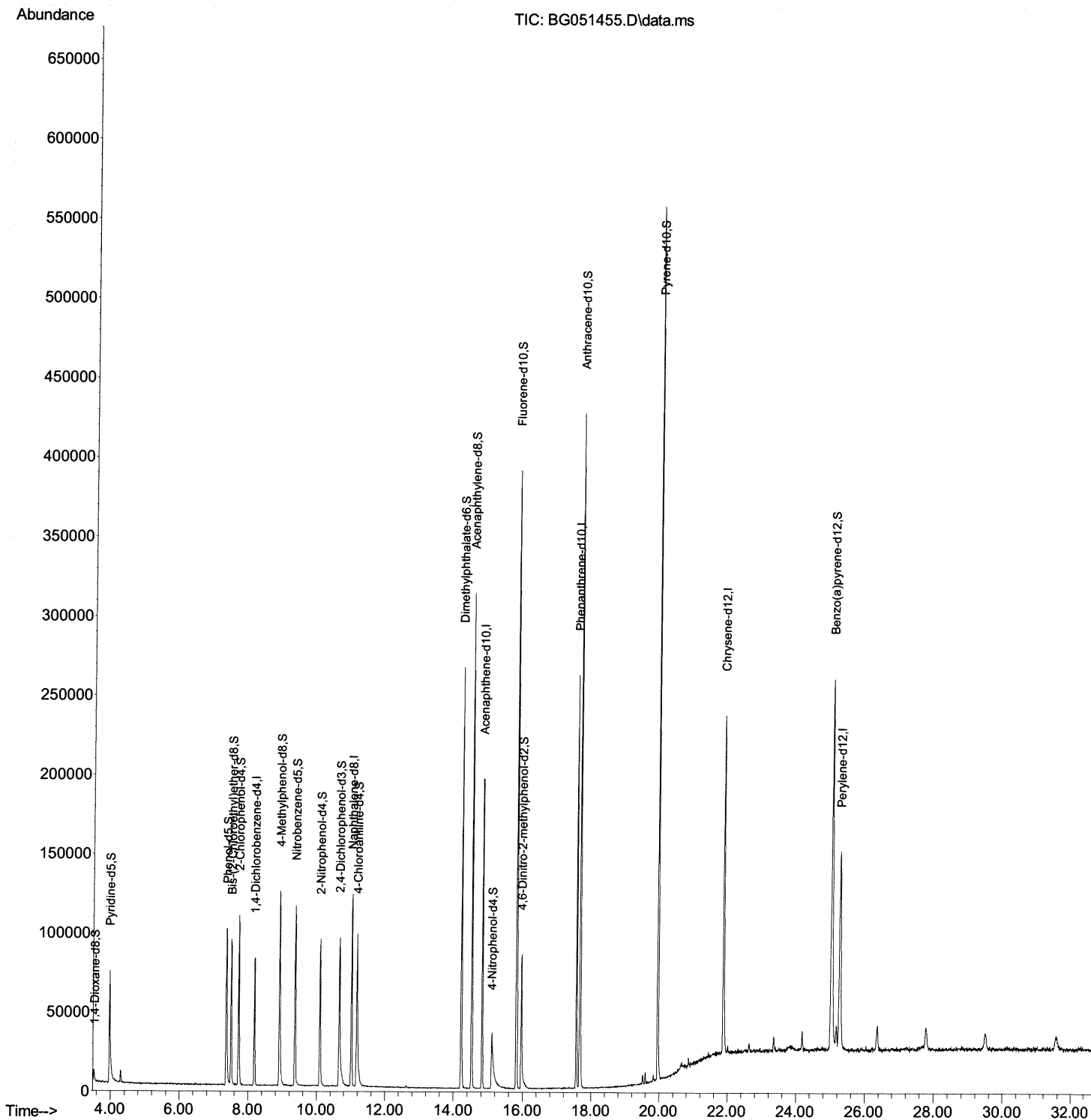
```
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120921\  
Data File : BG051455.D  
Acq On    : 10 Dec 2021  14:13  
Operator  : CG/JU  
Sample    : PB141295BL  
Misc      :  
ALS Vial  : 4    Sample Multiplier: 1
```

Instrument :
BNA_G
ClientSampleId :
SBLK295

Manual IntegrationsAPPROVED

Quant Time: Dec 11 01:29:23 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

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



Quantitation Report (Qedit)

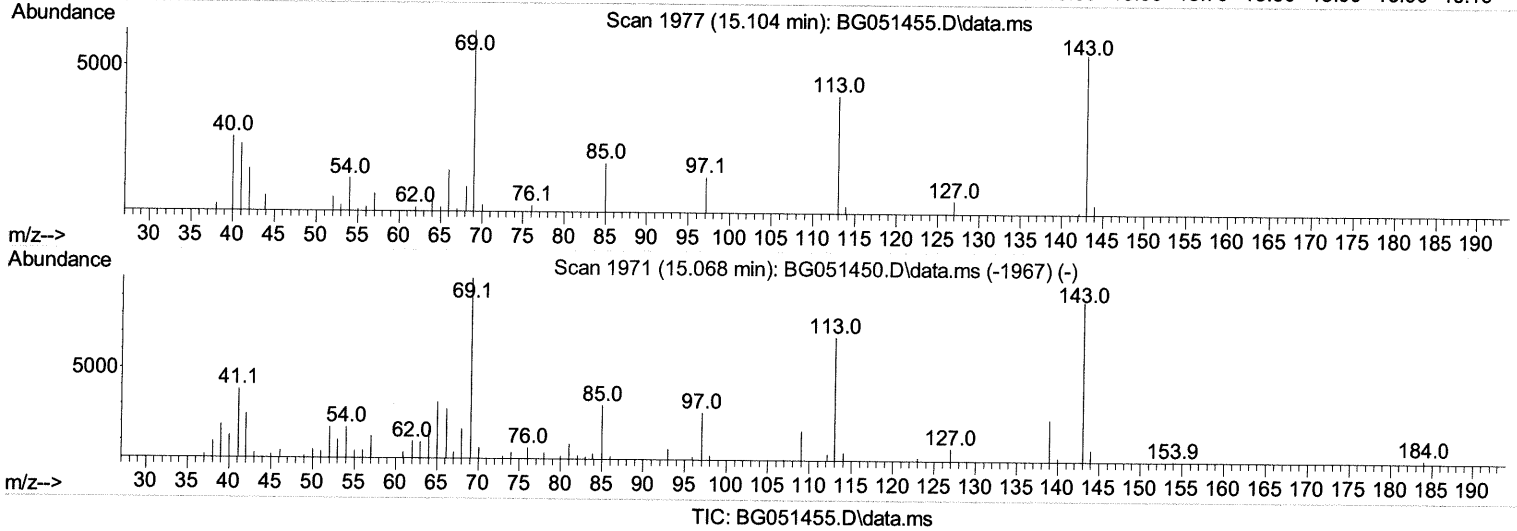
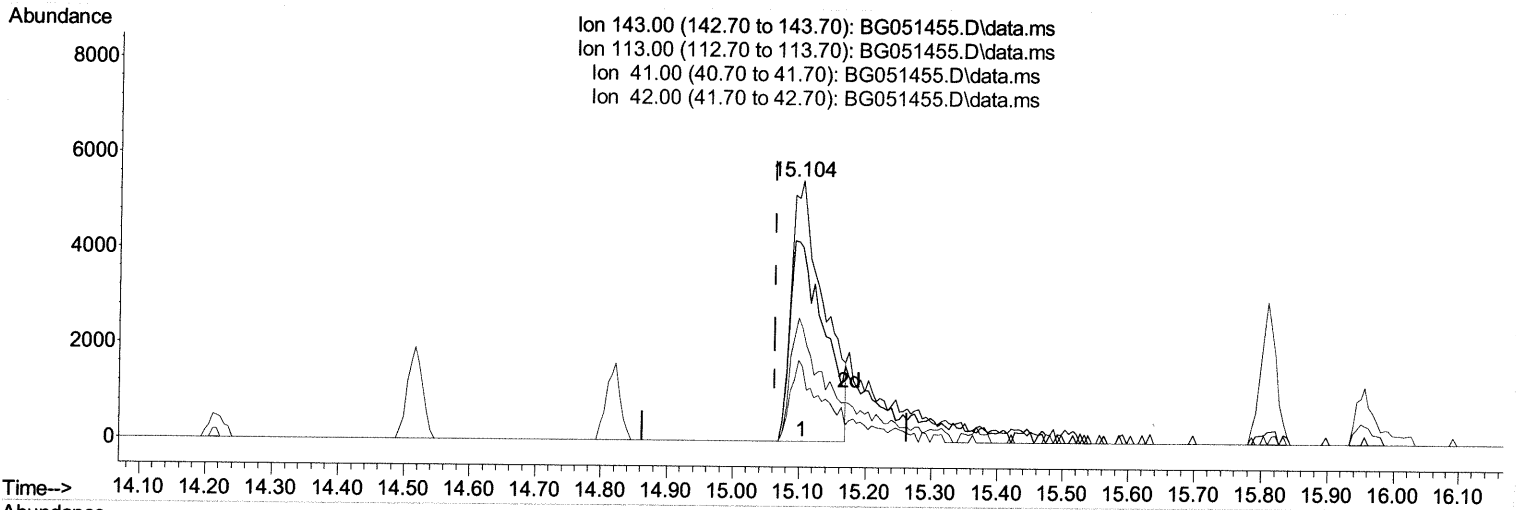
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(54) 4-Nitrophenol-d4 (S)

15.104min (+ 0.041) 21.94 ng/ul

response 18339

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	80.30	74.43
41.00	44.40	43.55
42.00	29.70	28.57

Quantitation Report (Qedit)

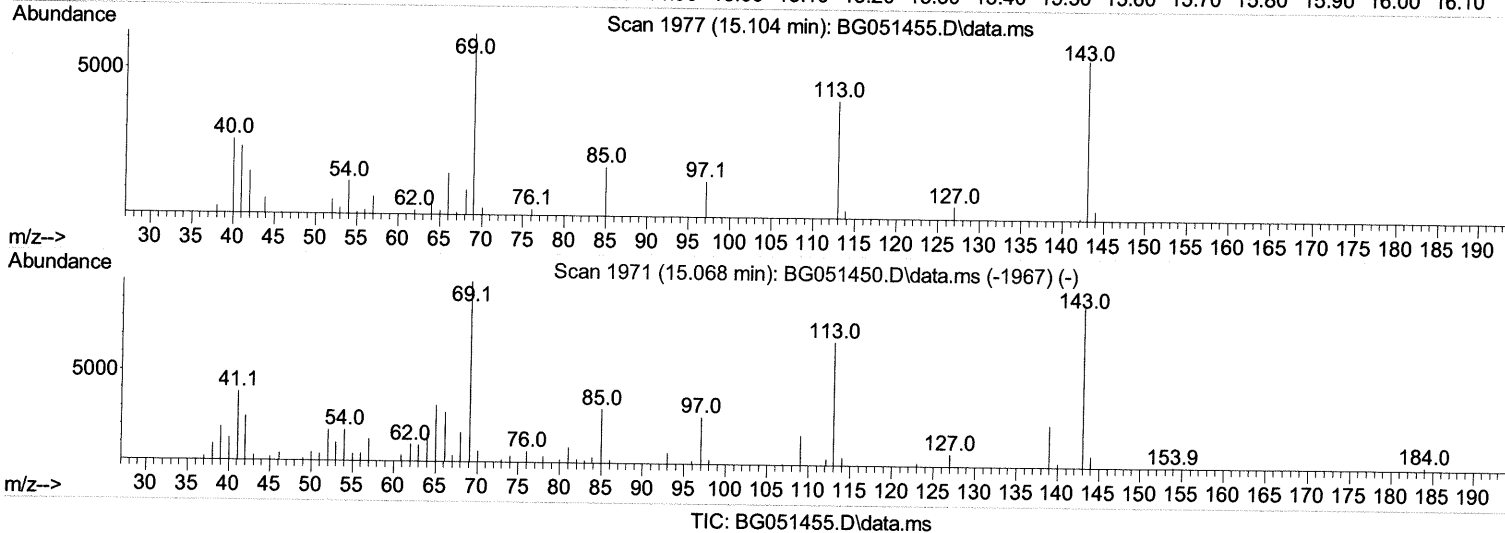
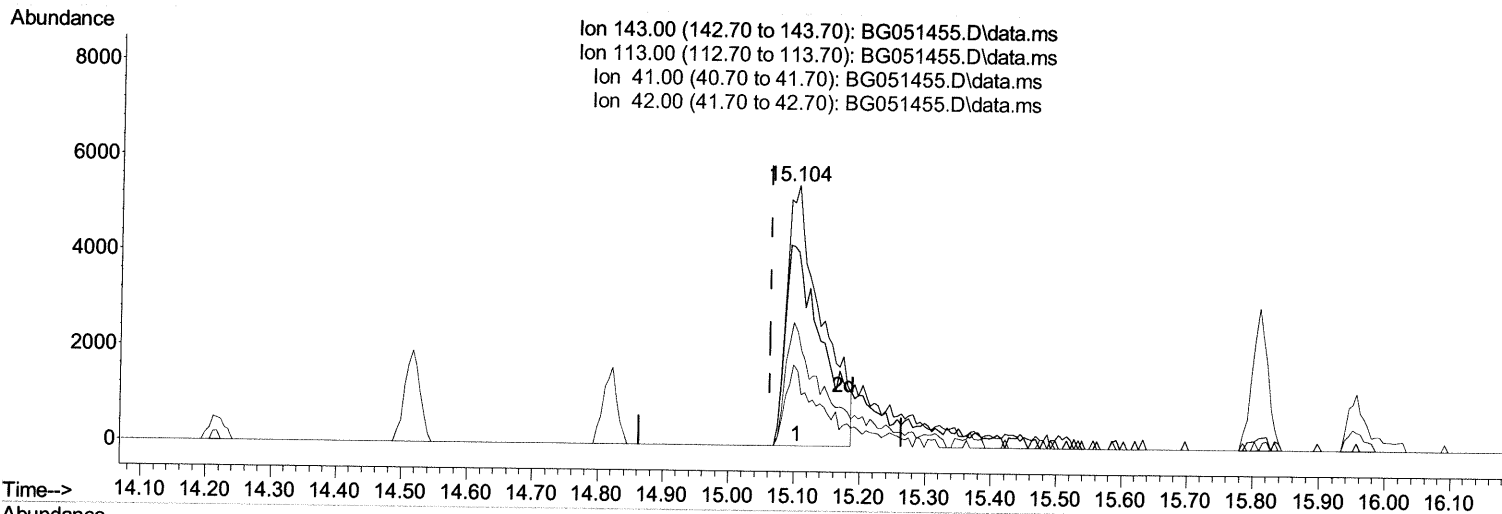
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 Data File : BG051455.D
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 Operator : CG/JU
 Sample : PB141295BL
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
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 ClientSampleId :
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 Supervised By :Yogesh Patel 12/15/2021



(54) 4-Nitrophenol-d4 (S)

15.104min (+ 0.041) 23.75 ng/ul m 12/11/21 JU

response 19847

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	80.30	74.43
41.00	44.40	43.55
42.00	29.70	28.57

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120921\
 Data File : BG051455.D
 Acq On : 10 Dec 2021 14:13
 Operator : CG/JU
 Sample : PB141295BL
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
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Manual IntegrationsAPPROVED

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.189	152	23962	20.000	ng/ul	0.00
20) Naphthalene-d8	11.015	136	105463	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.816	164	71741	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.572	188	168705	20.000	ng/ul	0.00
79) Chrysene-d12	21.873	240	155158	20.000	ng/ul	0.00
88) Perylene-d12	25.269	264	139044	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.535	96	4537	6.218	ng/uL	0.00
4) Pyridine-d5	3.970	84	52311	24.966	ng/ul	0.00
7) Phenol-d5	7.366	99	73929	30.306	ng/ul	0.01
9) Bis-(2-Chloroethyl)eth...	7.501	67	49427	31.596	ng/ul	0.00
11) 2-Chlorophenol-d4	7.724	132	53887	31.049	ng/ul	0.00
15) 4-Methylphenol-d8	8.917	113	54481	28.430	ng/ul	0.00
21) Nitrobenzene-d5	9.370	128	28970	31.666	ng/ul	0.00
24) 2-Nitrophenol-d4	10.098	143	31982	30.894	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.656	165	50044	29.715	ng/ul	0.00
31) 4-Chloroaniline-d4	11.173	131	75541	30.667	ng/ul	0.01
46) Dimethylphthalate-d6	14.217	166	188976	34.042	ng/ul	0.00
49) Acenaphthylene-d8	14.516	160	232676	33.094	ng/ul	0.00
54) 4-Nitrophenol-d4	15.104	143	19847m	23.747	ng/ul	0.04
60) Fluorene-d10	15.809	176	165787	33.549	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.956	200	24543	24.481	ng/ul	0.00
73) Anthracene-d10	17.672	188	273693	34.673	ng/ul	0.00
81) Pyrene-d10	19.951	212	326171	34.974	ng/ul	0.00
92) Benzo(a)pyrene-d12	25.034	264	267106	37.244	ng/ul	0.00

Target Compounds Qvalue

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