

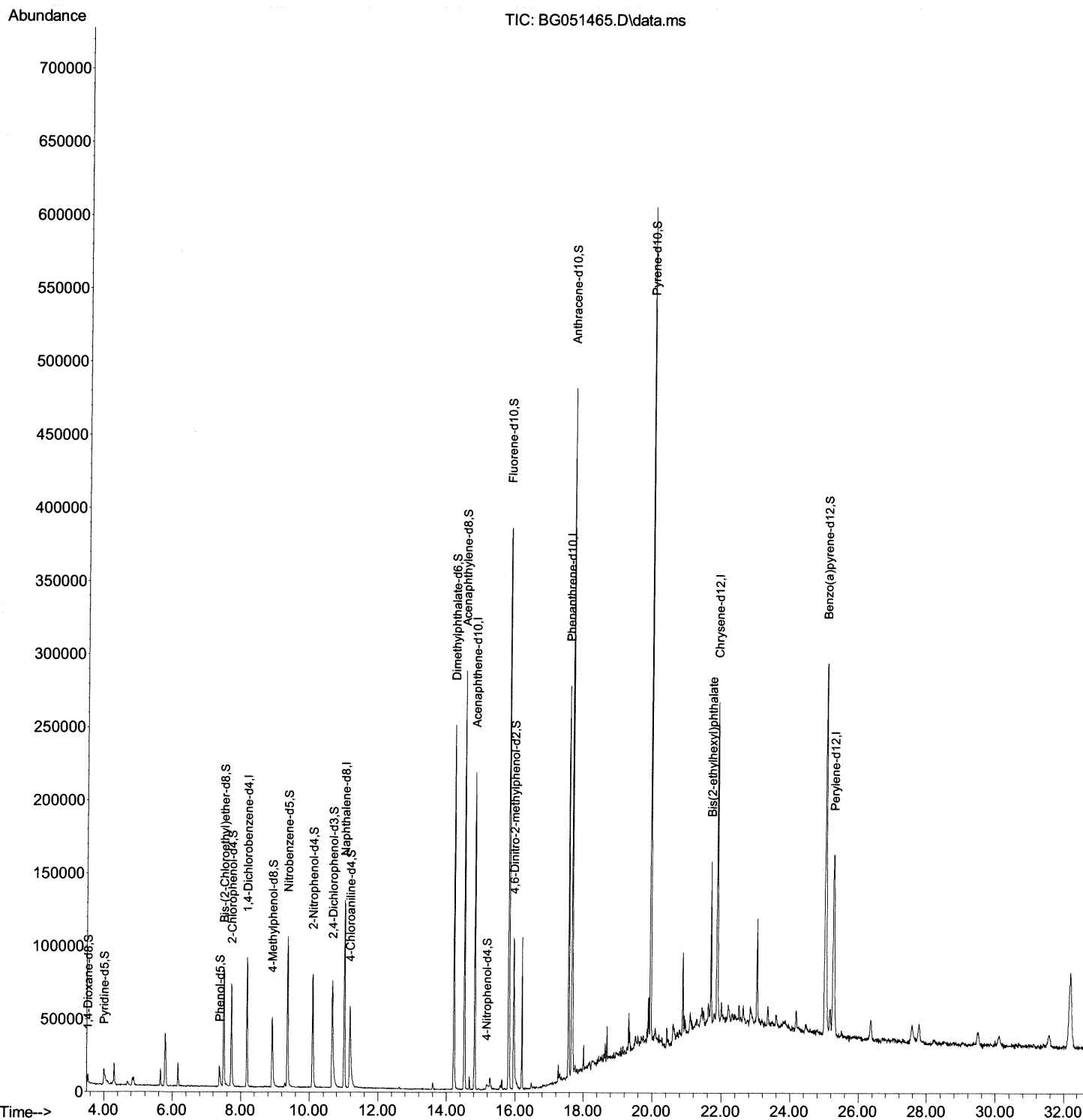
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120921\
Data File : BG051465.D
Acq On : 10 Dec 2021 21:03
Operator : CG/JU
Sample : M4985-12
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
EW5Q6

Manual IntegrationsAPPROVED

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

Quant Time: Dec 11 04:17:59 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)

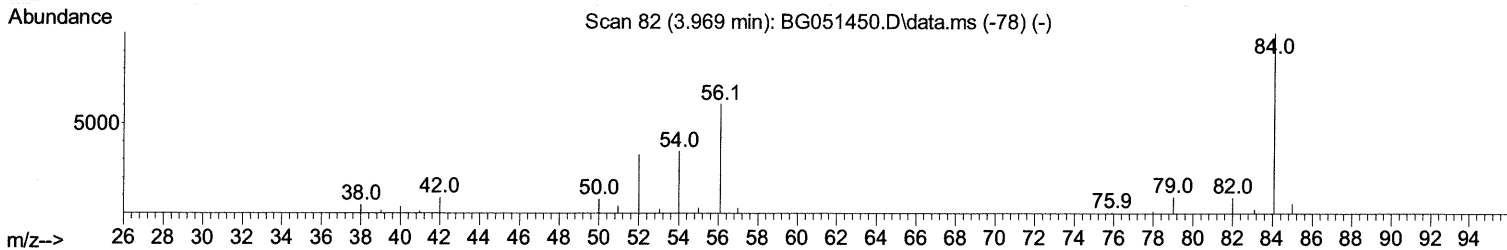
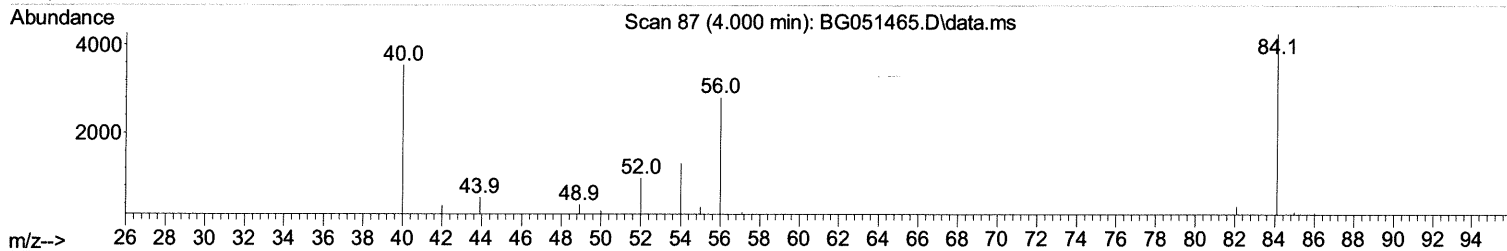
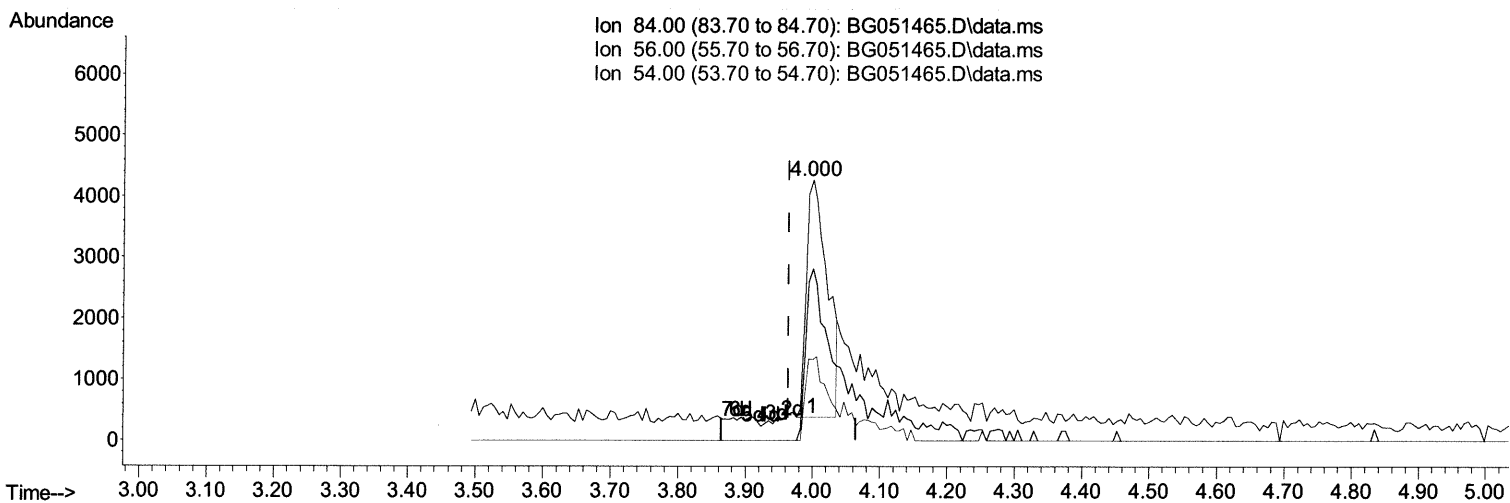
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TIC: BG051465.D\data.ms

(4) Pyridine-d5 (S)

4.000min (+ 0.035) 3.82 ng/ul

response 8474

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	68.00	65.86
54.00	31.50	30.93
0.00	0.00	0.00

Quantitation Report (Qedit)

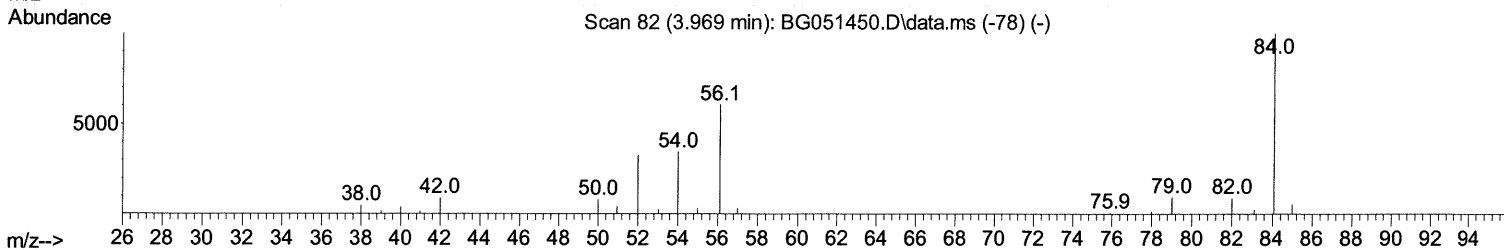
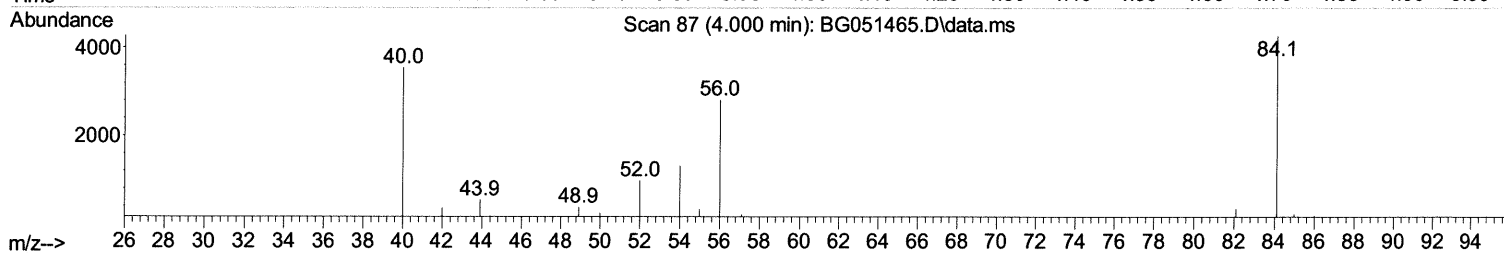
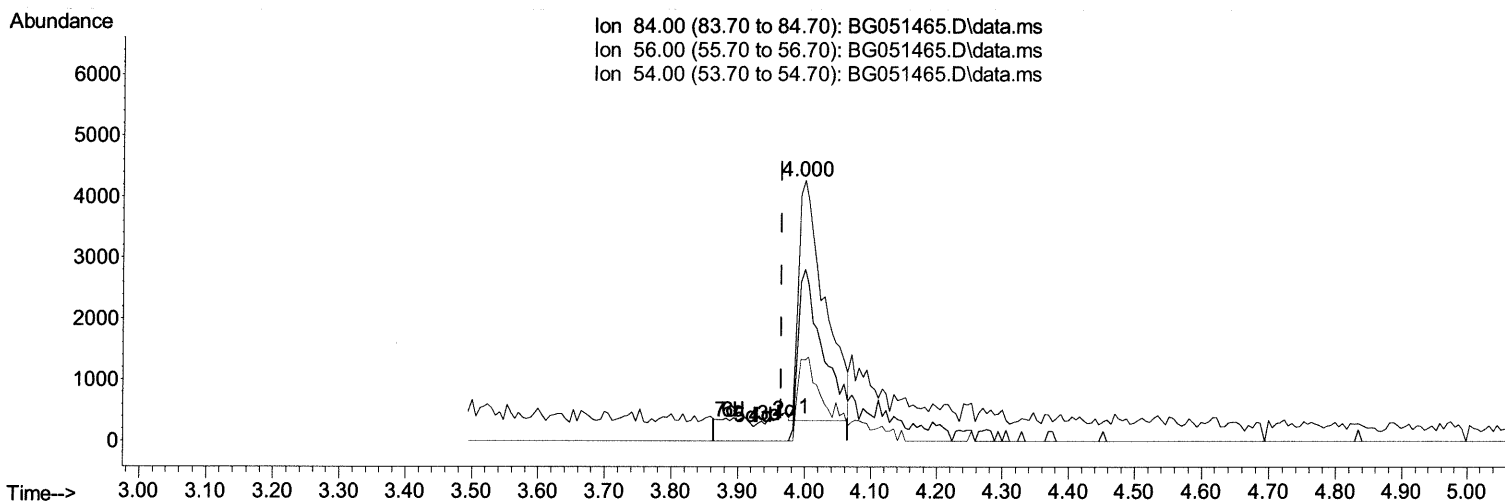
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TIC: BG051465.D\data.ms

(4) Pyridine-d5 (S)

4.000min (+ 0.035) 4.80 ng/ul m 12/16/21 ju

response 10631

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	68.00	65.86
54.00	31.50	30.93
0.00	0.00	0.00

Quantitation Report (Qedit)

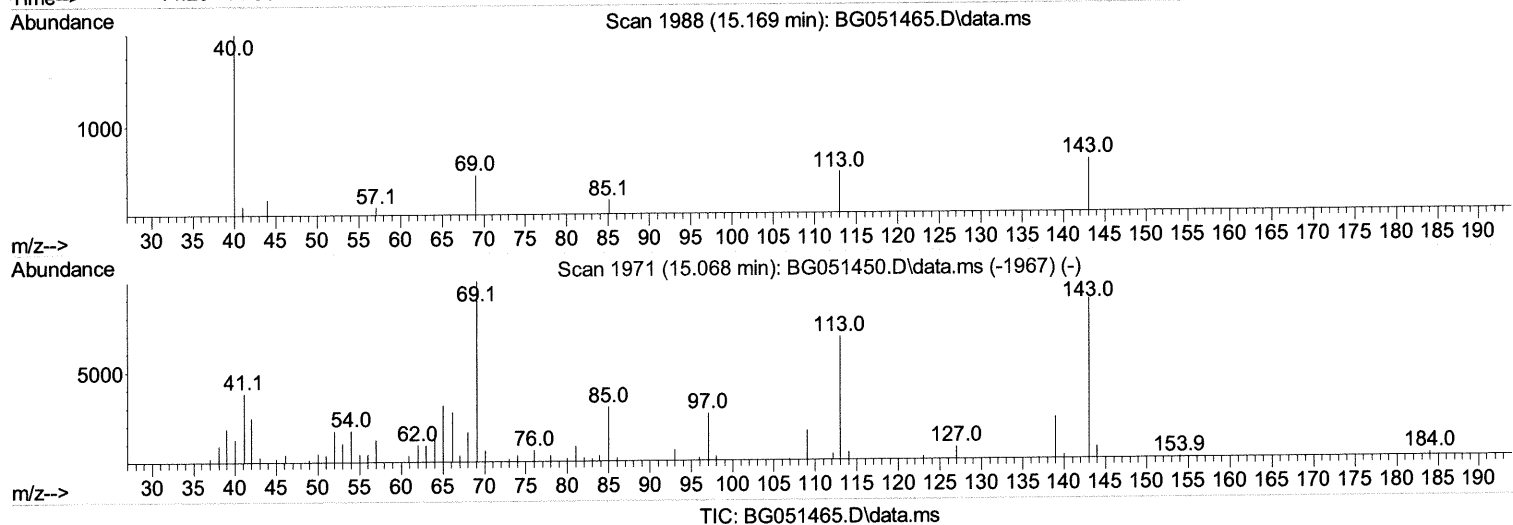
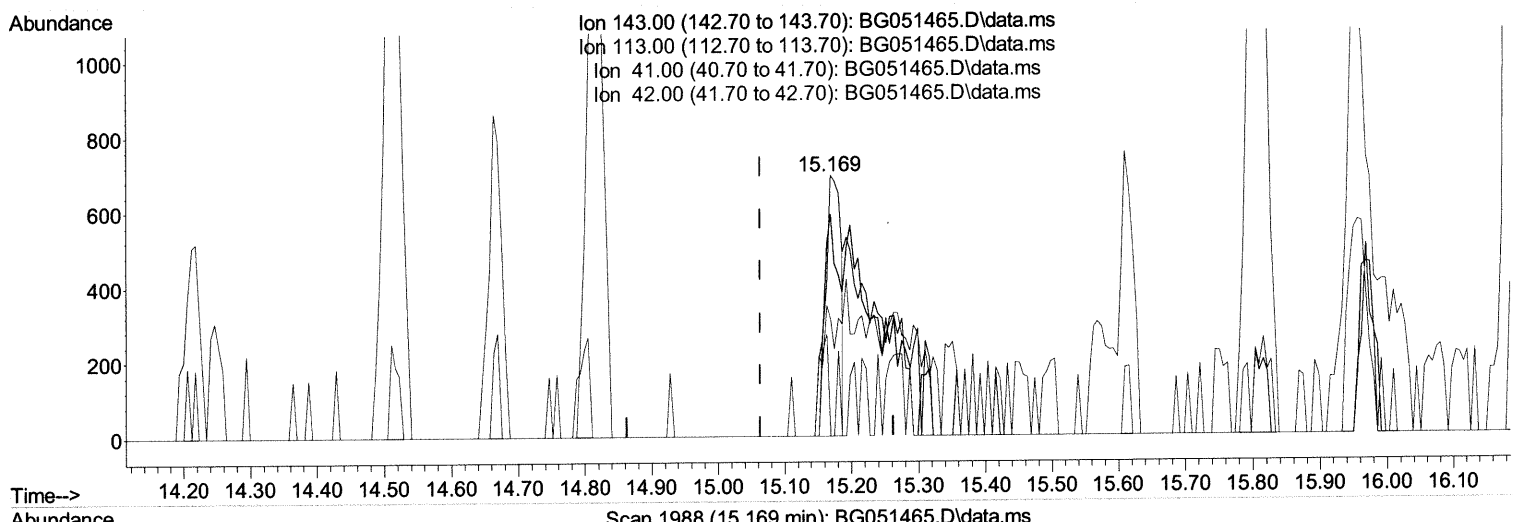
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(54) 4-Nitrophenol-d4 (S)

15.169min (+ 0.106) 1.27 ng/ul

response 1125

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	80.30	85.14
41.00	44.40	44.88
42.00	29.70	0.00#

Quantitation Report (Qedit)

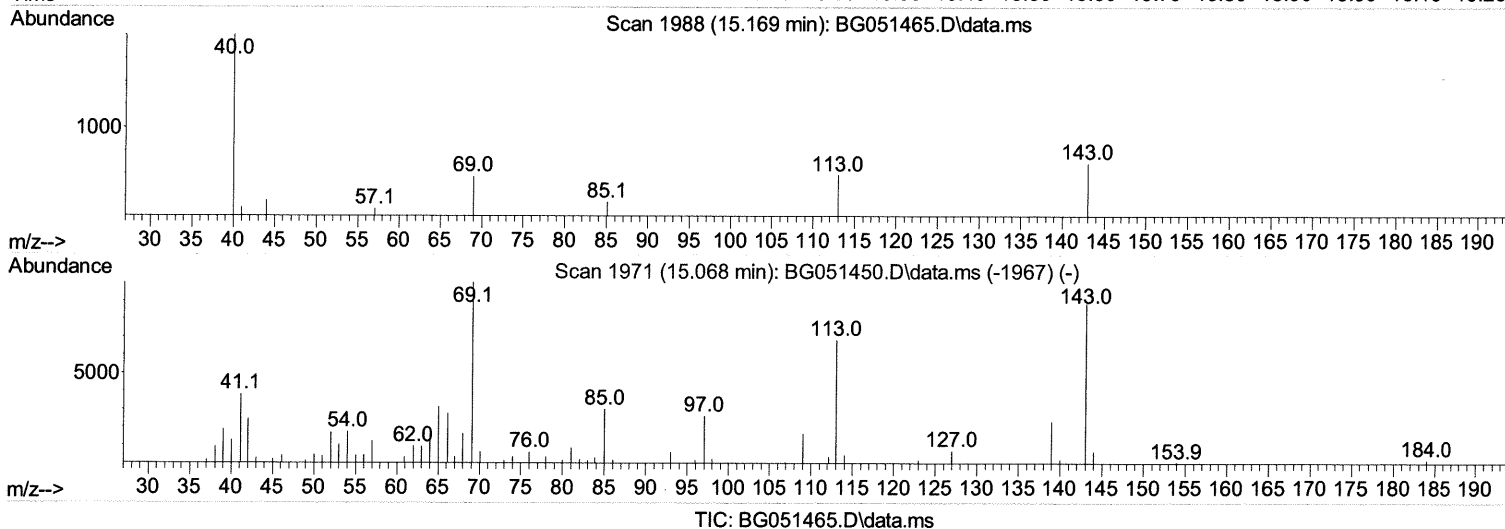
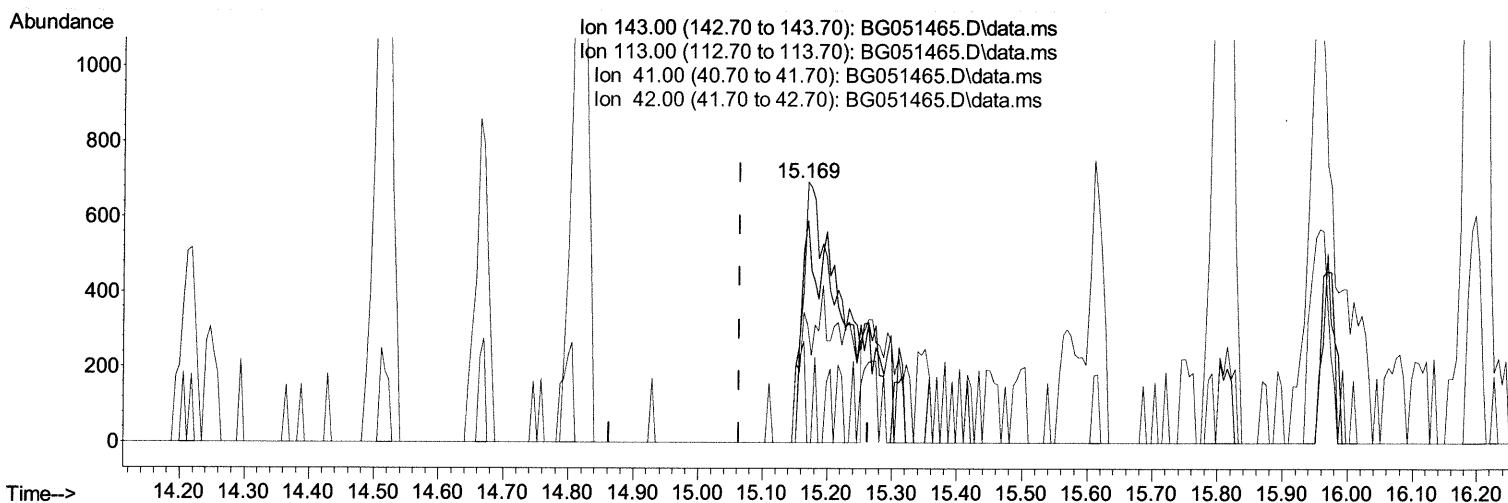
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 ALS Vial : 14 Sample Multiplier: 1

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



(54) 4-Nitrophenol-d4 (S)

15.169min (+ 0.106) 2.92 ng/ul m 12/16/21 JU

response 2577

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	80.30	85.14
41.00	44.40	44.88
42.00	29.70	0.00#

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 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 EW5Q6

Manual Integrations APPROVED

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.183	152	25344	20.000	ng/ul	0.00
20) Naphthalene-d8	11.009	136	110981	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.816	164	75798	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.566	188	168800	20.000	ng/ul	0.00
79) Chrysene-d12	21.867	240	150621	20.000	ng/ul	0.00
88) Perylene-d12	25.263	264	144183	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.530	96	3549	4.599	ng/ul	0.00
4) Pyridine-d5	4.000	84	10631m >	4.797	ng/ul >	0.04 2/1/21 JU
7) Phenol-d5	7.384	99	12806	4.963	ng/ul	0.03
9) Bis-(2-Chloroethyl)eth...	7.501	67	45591	27.555	ng/ul	0.00
11) 2-Chlorophenol-d4	7.731	132	38881	21.181	ng/ul	0.00
15) 4-Methylphenol-d8	8.923	113	26081	12.868	ng/ul	0.01
21) Nitrobenzene-d5	9.370	128	26763	27.799	ng/ul	0.00
24) 2-Nitrophenol-d4	10.098	143	28680	26.326	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.662	165	41108	23.195	ng/ul	0.01
31) 4-Chloroaniline-d4	11.174	131	49912	19.255	ng/ul	0.01
46) Dimethylphthalate-d6	14.211	166	181512	30.947	ng/ul	0.00
49) Acenaphthylene-d8	14.517	160	223451	30.081	ng/ul	0.00
54) 4-Nitrophenol-d4	15.169	143	2577m >	2.918	ng/ul >	0.11 2/1/21 JU
60) Fluorene-d10	15.809	176	166181	31.829	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.950	200	25867	25.788	ng/ul	0.00
73) Anthracene-d10	17.666	188	281401	35.629	ng/ul	0.00
81) Pyrene-d10	19.946	212	322797	35.655	ng/ul	0.00
92) Benzo(a)pyrene-d12	25.028	264	280415	37.707	ng/ul	0.00
Target Compounds						
86) Bis(2-ethylhexyl)phtha...	21.696	149	41212	6.219	ng/ul	Qvalue 100

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