

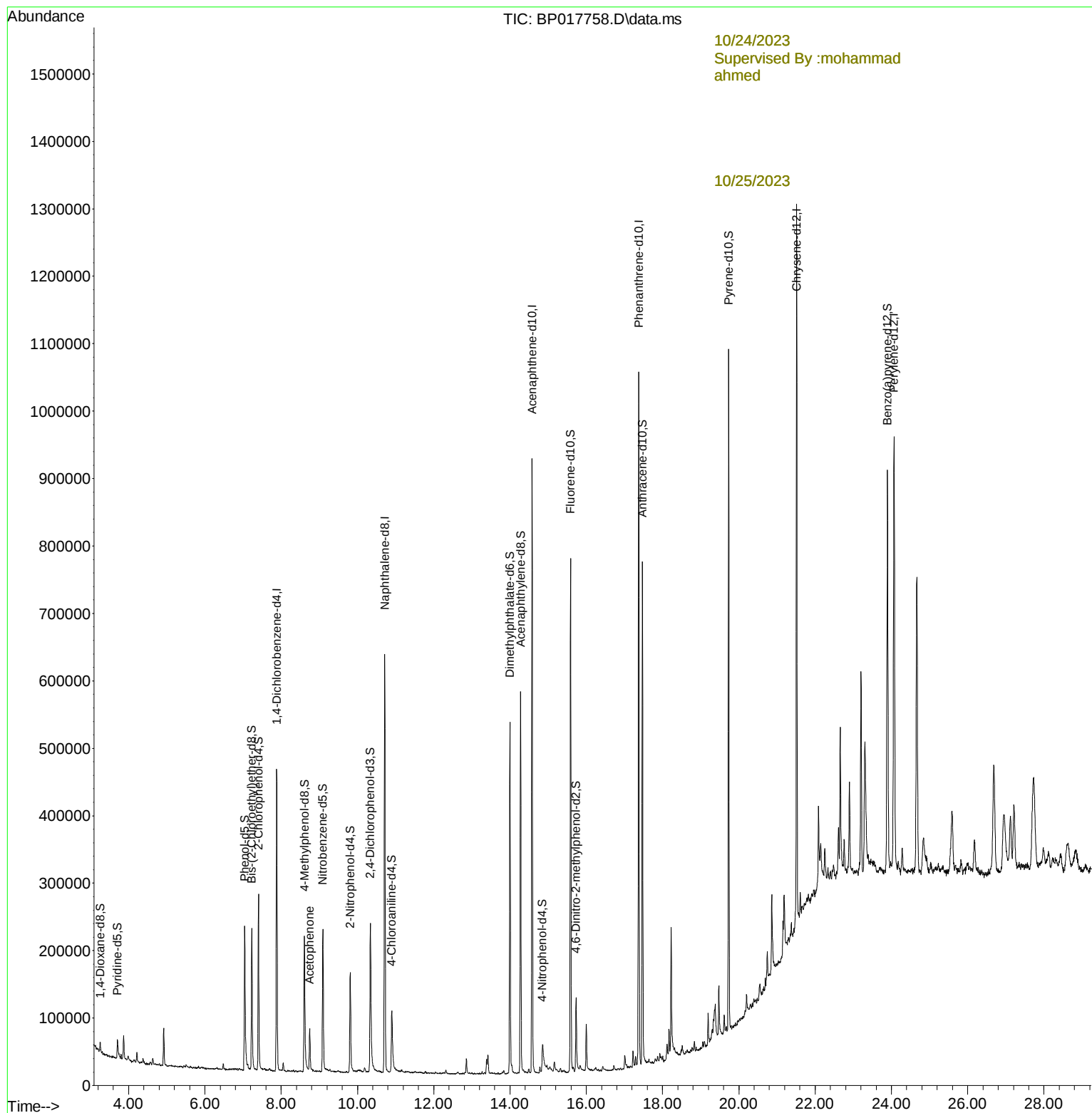
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102323\
Data File : BP017758.D
Acq On : 23 Oct 2023 20:01
Operator : MA/JU
Sample : 04926-15
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
GCD00

Manual IntegrationsAPPROVED

Quant Time: Oct 24 05:58:15 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP101823.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Thu Oct 19 01:32:15 2023
Response via : Initial Calibration

Reviewed By :Yogesh Patel 10/24/2023
Supervised By :mohammad ahmed 10/25/2023



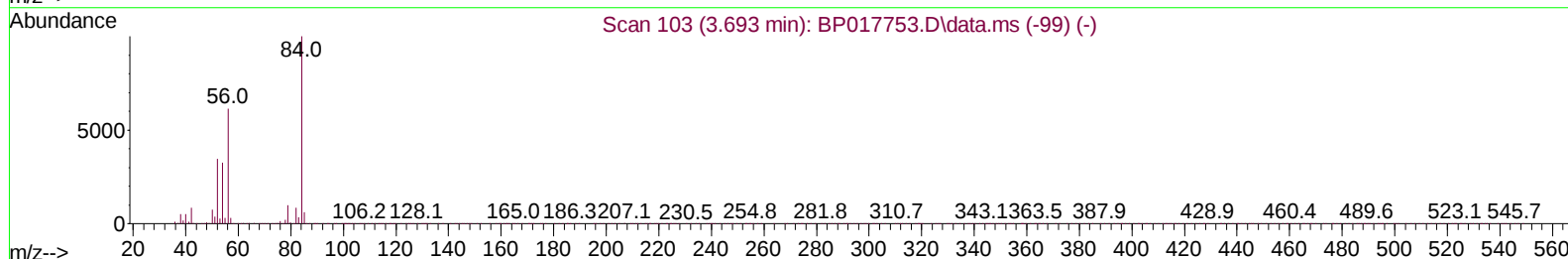
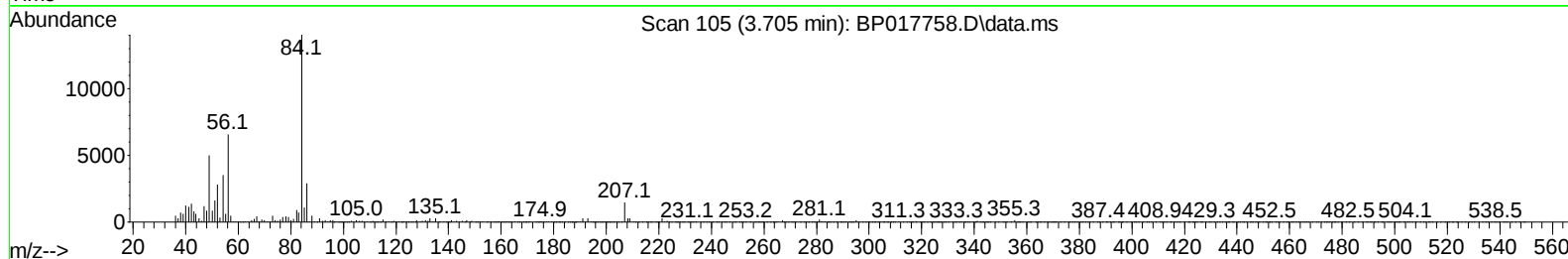
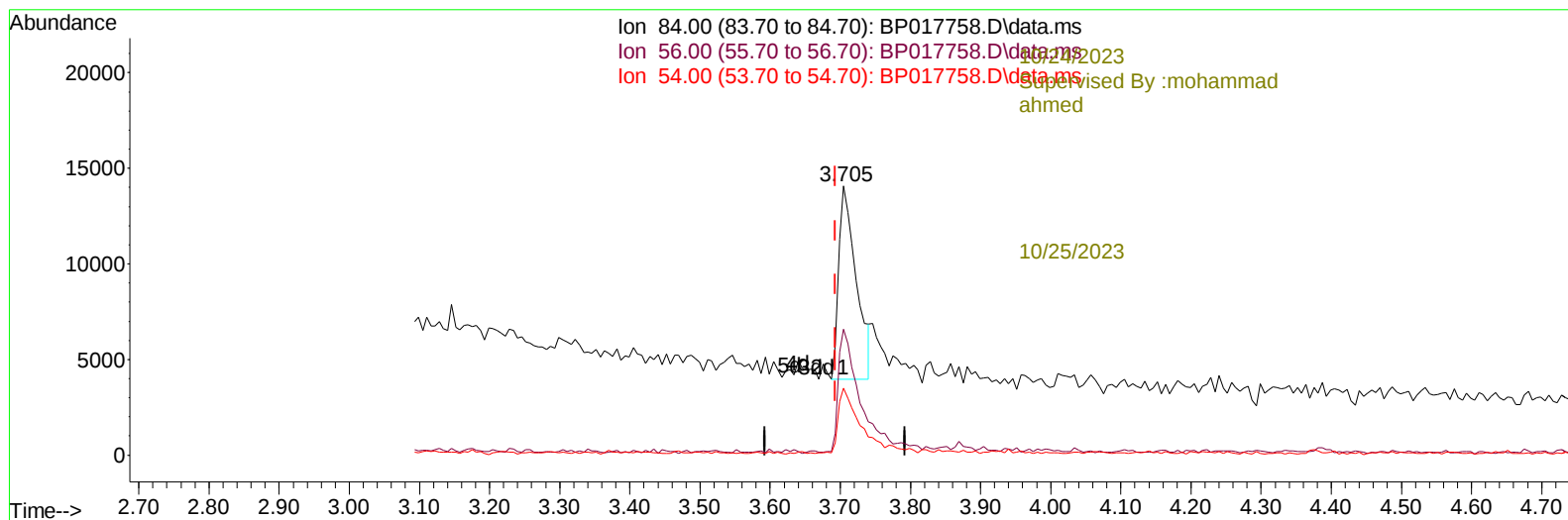
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TIC: BP017758.D\data.ms

(4) Pyridine-d5 (S)

3.705min (+ 0.012) 2.14 ng/u1

response 17660

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	55.10	46.83
54.00	30.40	24.94
0.00	0.00	0.00

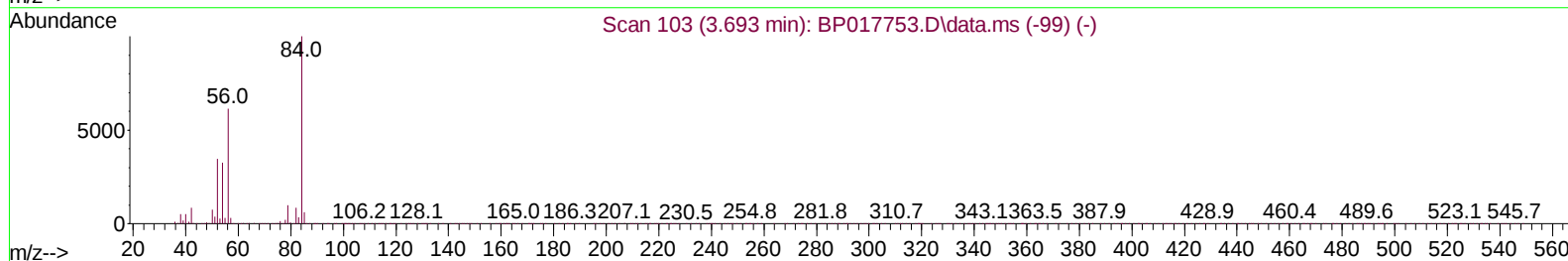
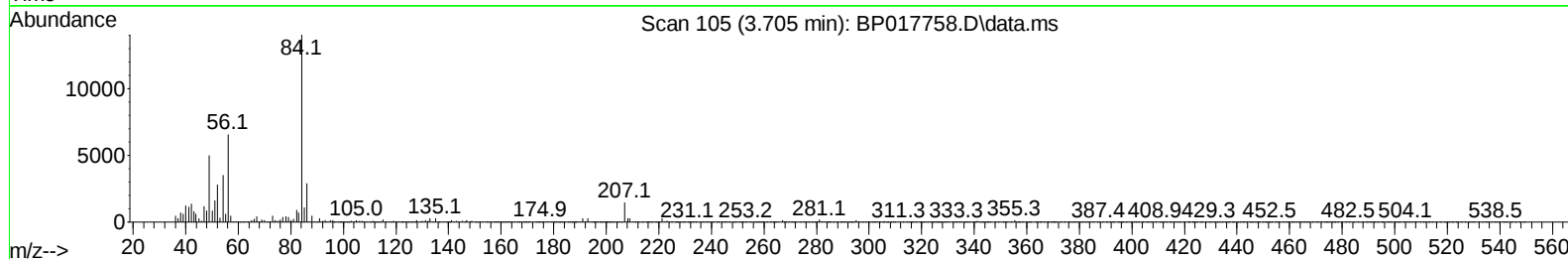
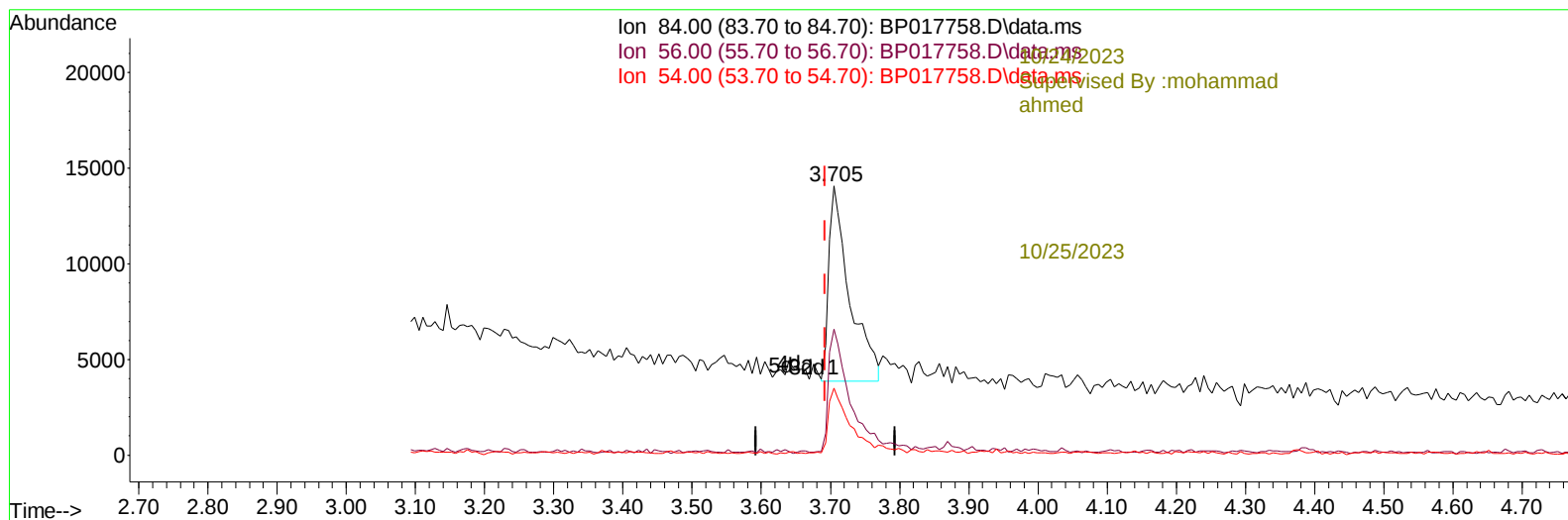
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TIC: BP017758.D\data.ms

(4) Pyridine-d5 (S)

3.705min (+ 0.012) 2.58 ng/u1 m

response 21265

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	55.10	46.83
54.00	30.40	24.94
0.00	0.00	0.00

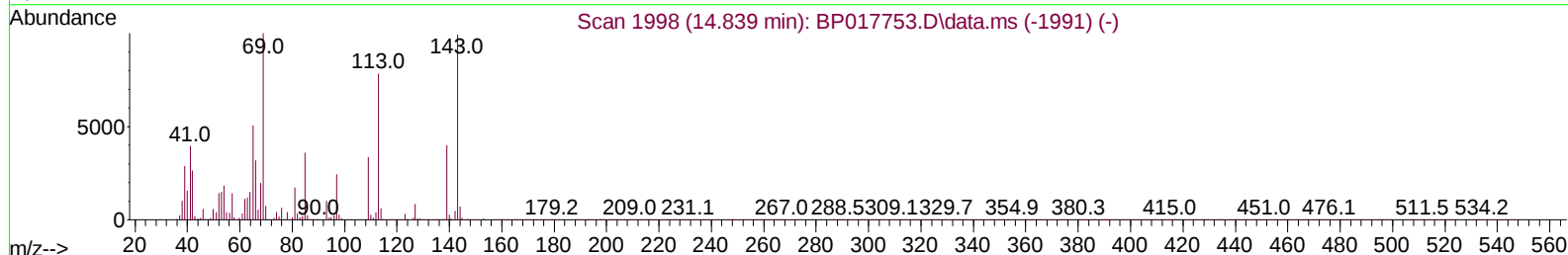
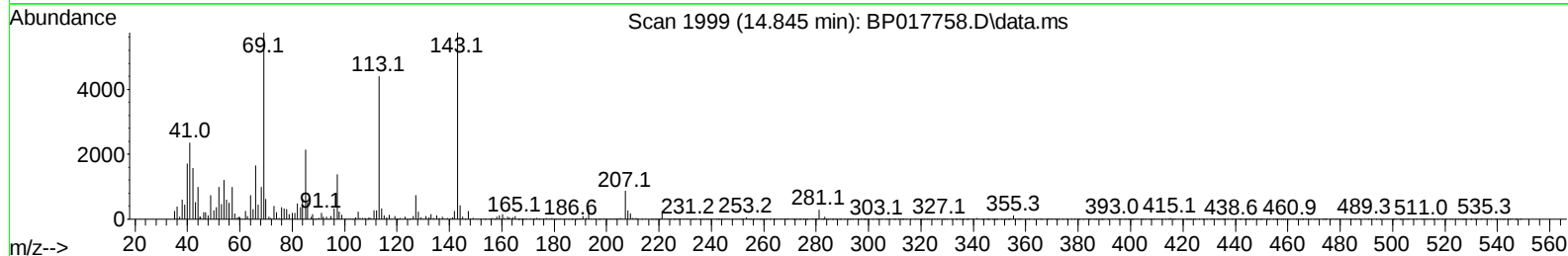
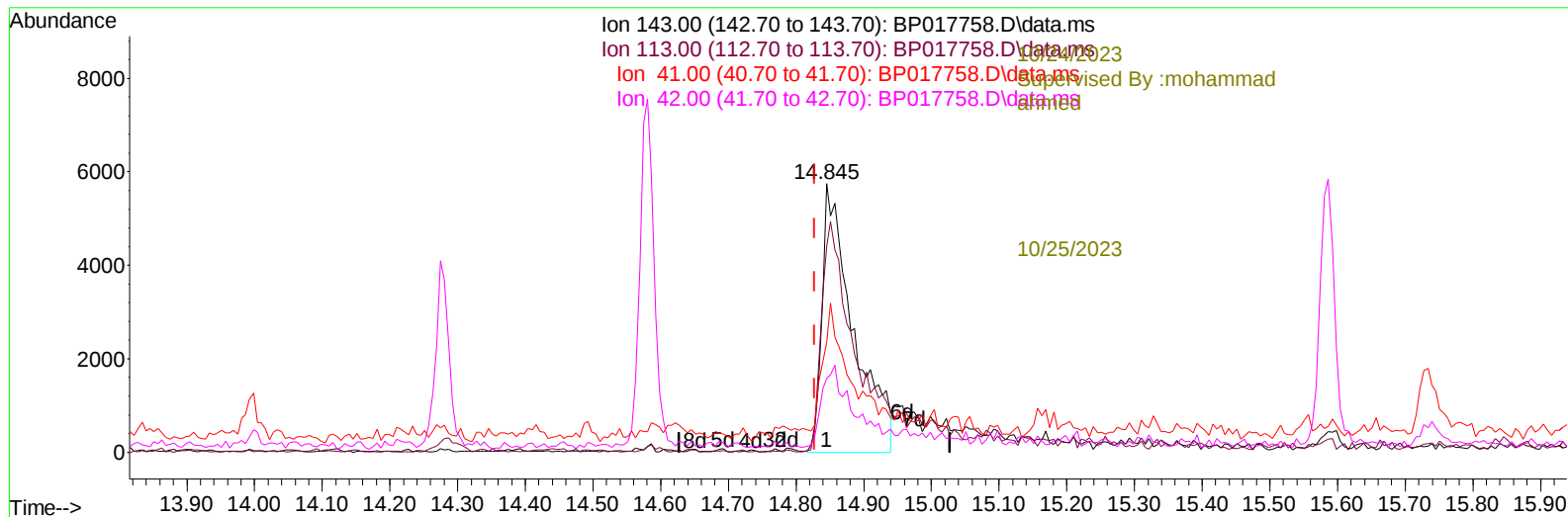
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TIC: BP017758.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.845min (+ 0.018) 5.39 ng/ul m

response 18267

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	81.70	76.71
41.00	39.20	41.03
42.00	25.60	27.39

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Compound	R.T.	Q	Ion	Response	Conc	Units	Dev(Min)
-----10/24/2023							
Supervised By :mohammad							
ahmed							
Internal Standards							
1) 1,4-Dichlorobenzene-d4	7.881	152		117051	20.000	ng/ul	0.00
20) Naphthalene-d8	10.716	136		468045	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.581	164		283105	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.375	188		588665	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.516	240		470578	20.000	ng/ul	0.00
88) Perylene-d12	24.068	264		474087	20.000	ng/ul	0.0025/2023
System Monitoring Compounds							
3) 1,4-Dioxane-d8	3.258	96		8421	2.676	ng/uL	0.00
4) Pyridine-d5	3.705	84		21265m	2.581	ng/ul	0.01
7) Phenol-d5	7.046	99		127490	13.089	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.228	67		93027	15.465	ng/ul	0.00
11) 2-Chlorophenol-d4	7.405	132		111789	14.071	ng/ul	0.00
15) 4-Methylphenol-d8	8.610	113		83088	10.862	ng/ul	0.00
21) Nitrobenzene-d5	9.093	128		47055	12.407	ng/ul	0.00
24) 2-Nitrophenol-d4	9.810	143		44389	10.593	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.340	165		94941	11.550	ng/ul	0.00
31) 4-Chloroaniline-d4	10.898	131		63794	5.930	ng/ul	0.00
46) Dimethylphthalate-d6	13.998	166		326753	13.749	ng/ul	0.00
49) Acenaphthylene-d8	14.281	160		359306	13.557	ng/ul	0.00
54) 4-Nitrophenol-d4	14.845	143		18267m	5.394	ng/ul	0.02
60) Fluorene-d10	15.586	176		281427	13.699	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.734	200		22725	5.861	ng/ul	0.00
73) Anthracene-d10	17.475	188		406767	13.522	ng/ul	0.00
81) Pyrene-d10	19.733	212		491560	16.661	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.898	264		381149	14.435	ng/ul	0.00
Target Compounds							
16) Acetophenone	8.746	105		32327	2.688	ng/ul	Qvalue 93

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