

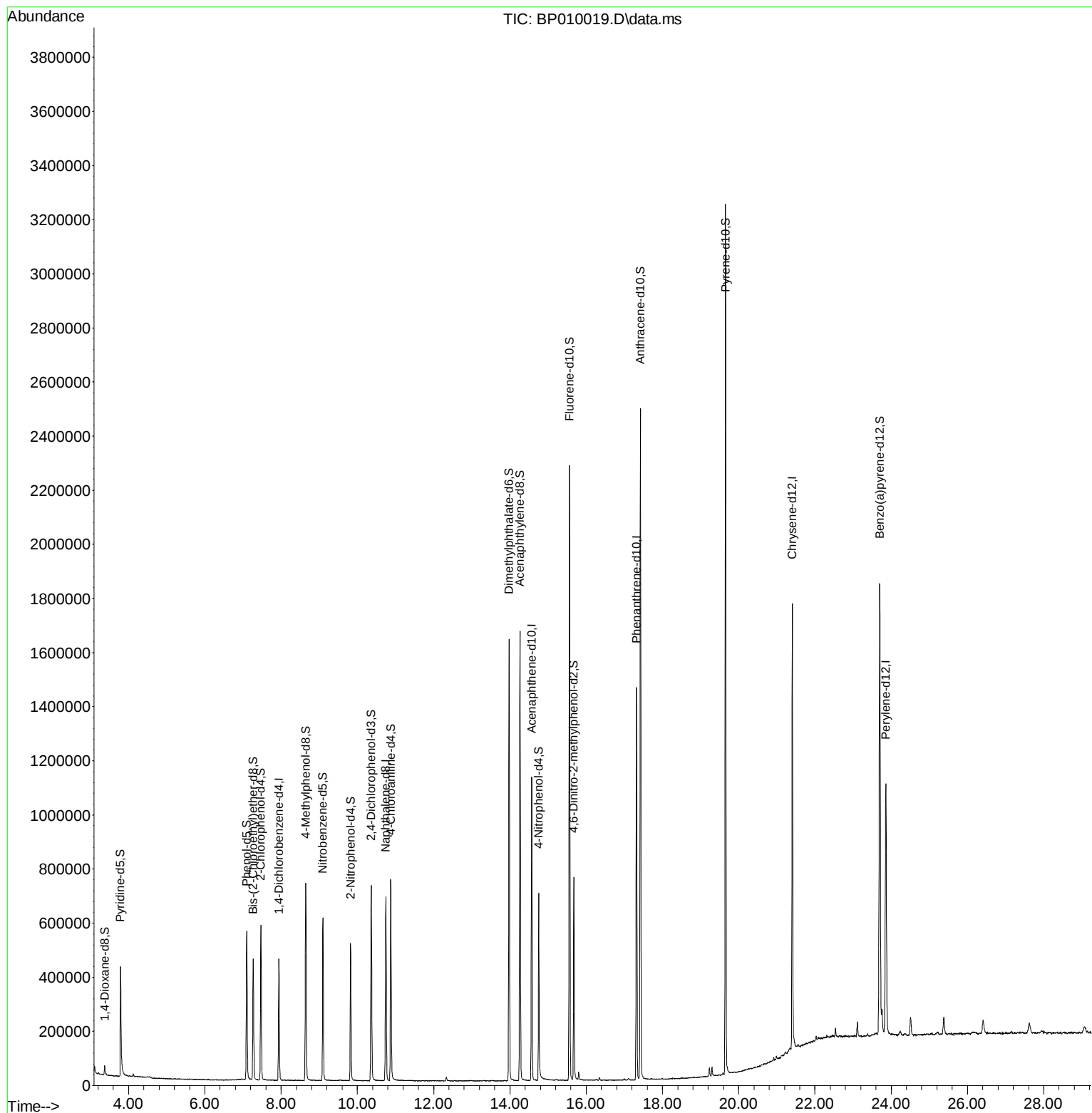
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042122\
Data File : BP010019.D
Acq On : 21 Apr 2022 17:53
Operator : CG/JU
Sample : PB144260BL
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
SBLK260

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 04/22/2022
Supervised By :mohammad ahmed 04/26/2022

Quant Time: Apr 22 04:18:29 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP042122.M
Quant Title : SVOA CALIBRATION
QLast Update : Thu Apr 21 14:49:38 2022
Response via : Initial Calibration



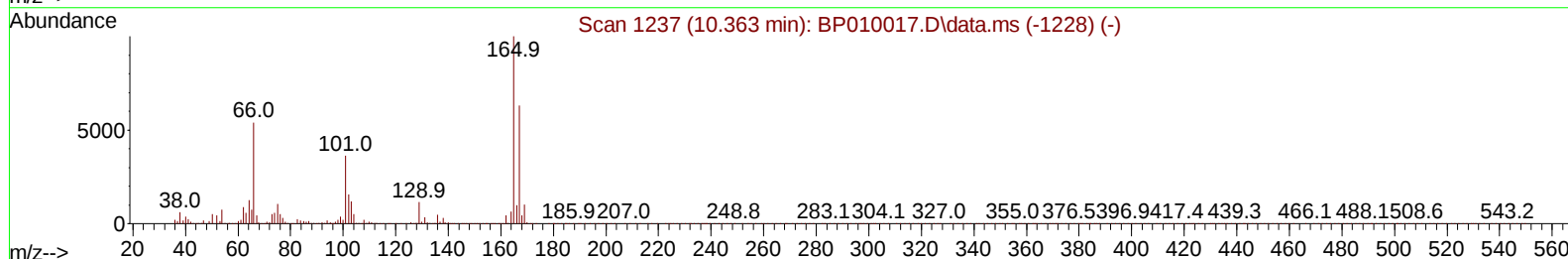
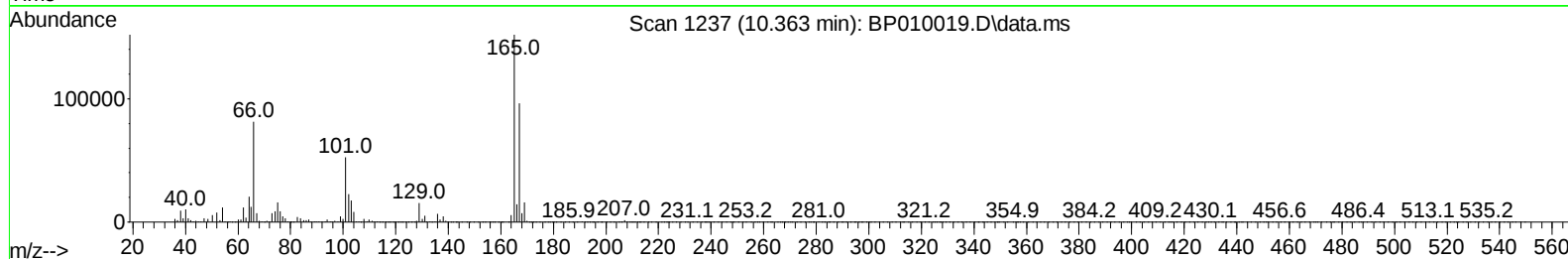
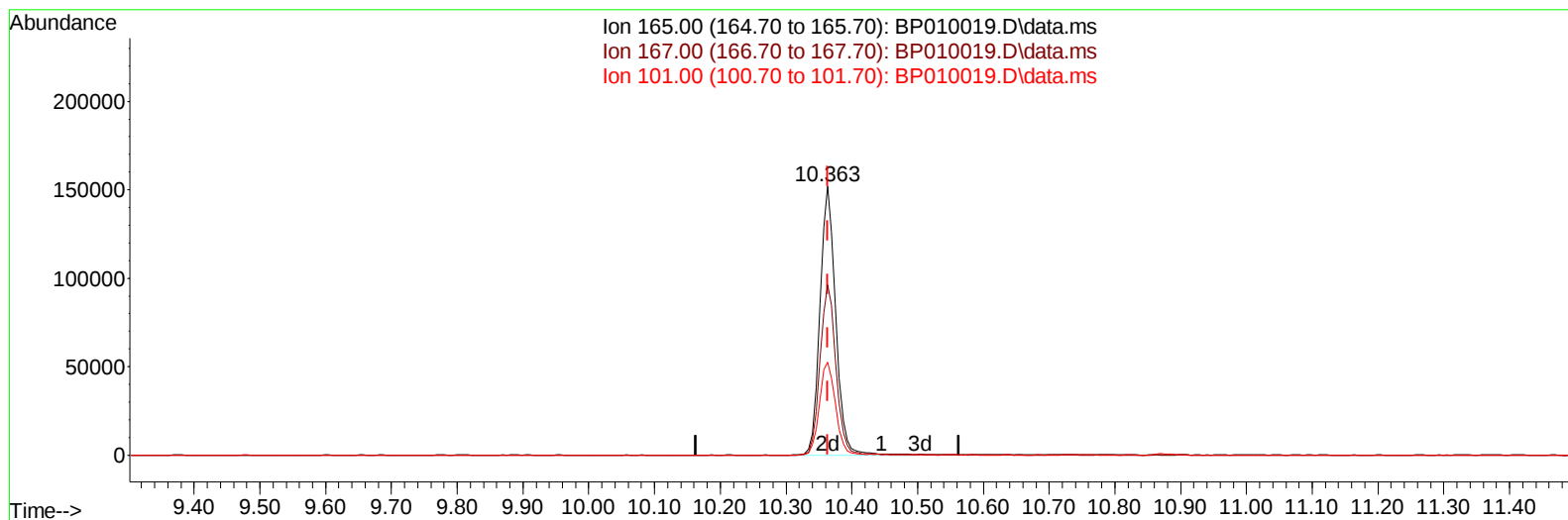
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TIC: BP010019.D\data.ms

(28) 2,4-Dichlorophenol-d3 (S)

10.363min (0.000) 28.72 ng/ul m

response 255052

Ion	Exp%	Act%
165.00	100.00	100.00
167.00	57.70	63.26
101.00	49.80	34.49#
0.00	0.00	0.00

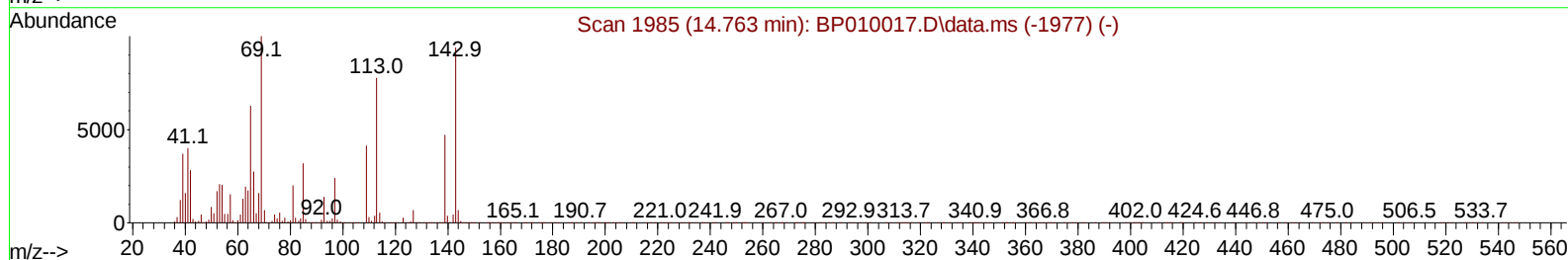
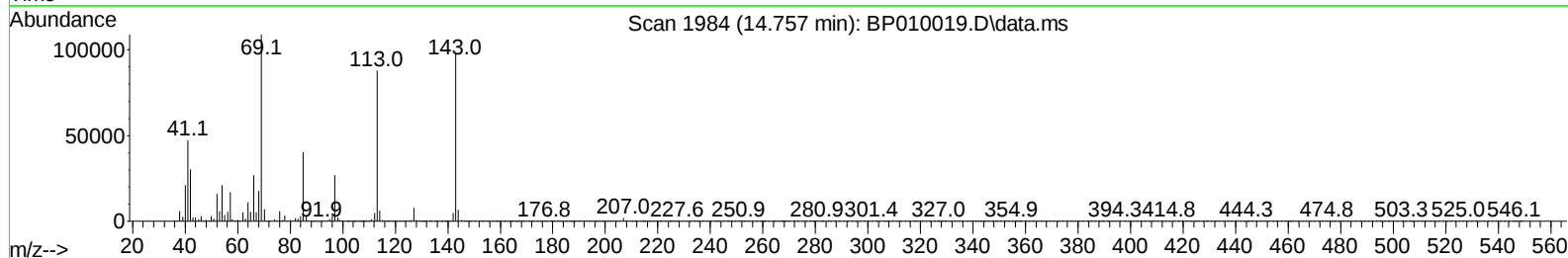
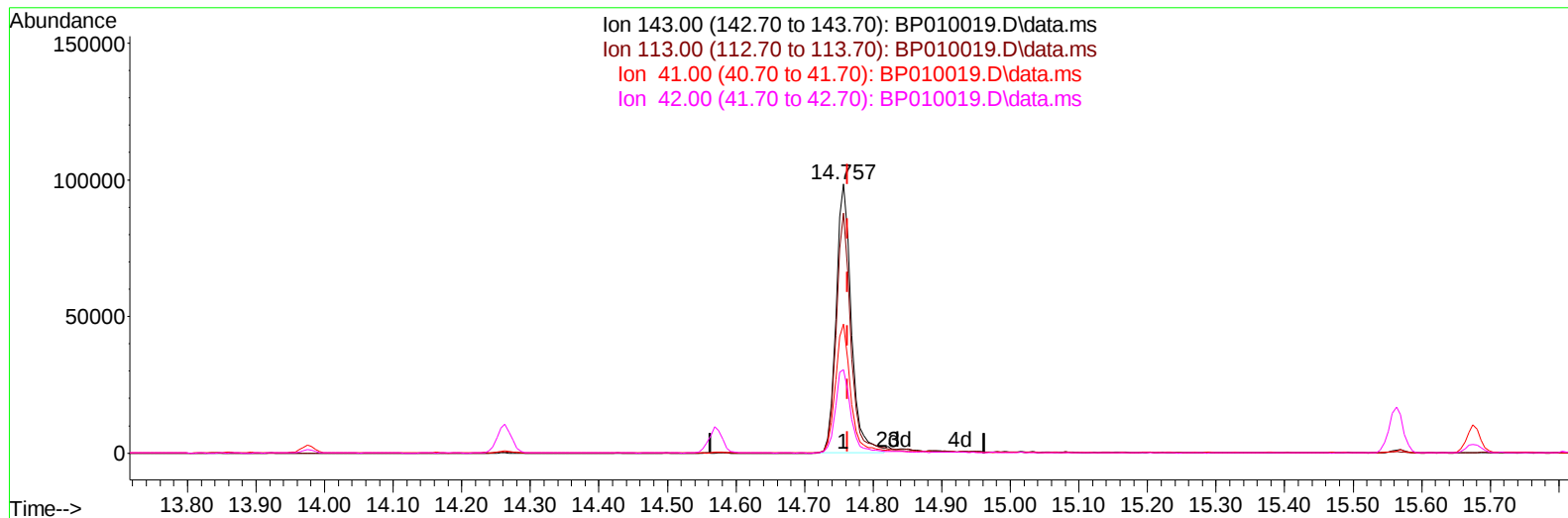
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TIC: BP010019.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.757min (-0.006) 26.74 ng/u1

response 151092

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	132.70	89.21#
41.00	23.20	48.07#
42.00	18.00	30.93#

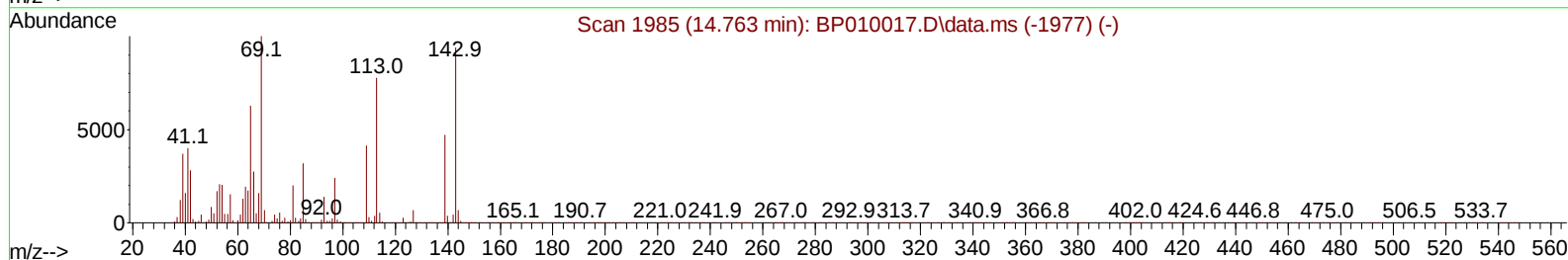
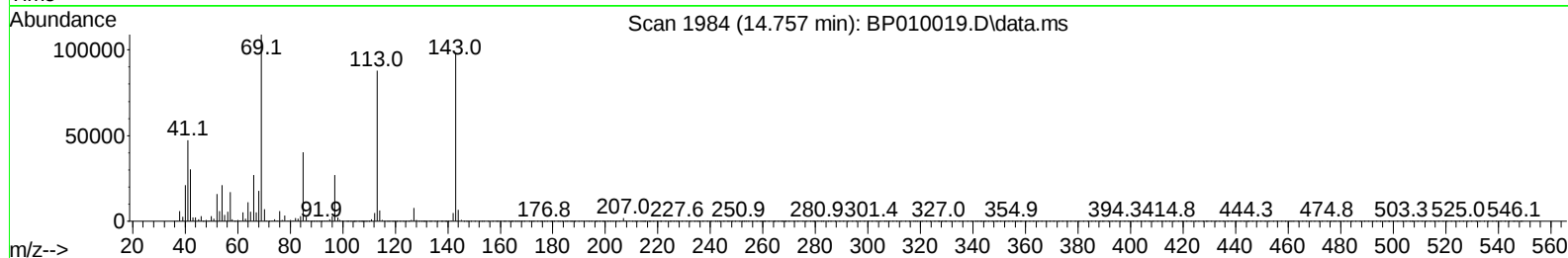
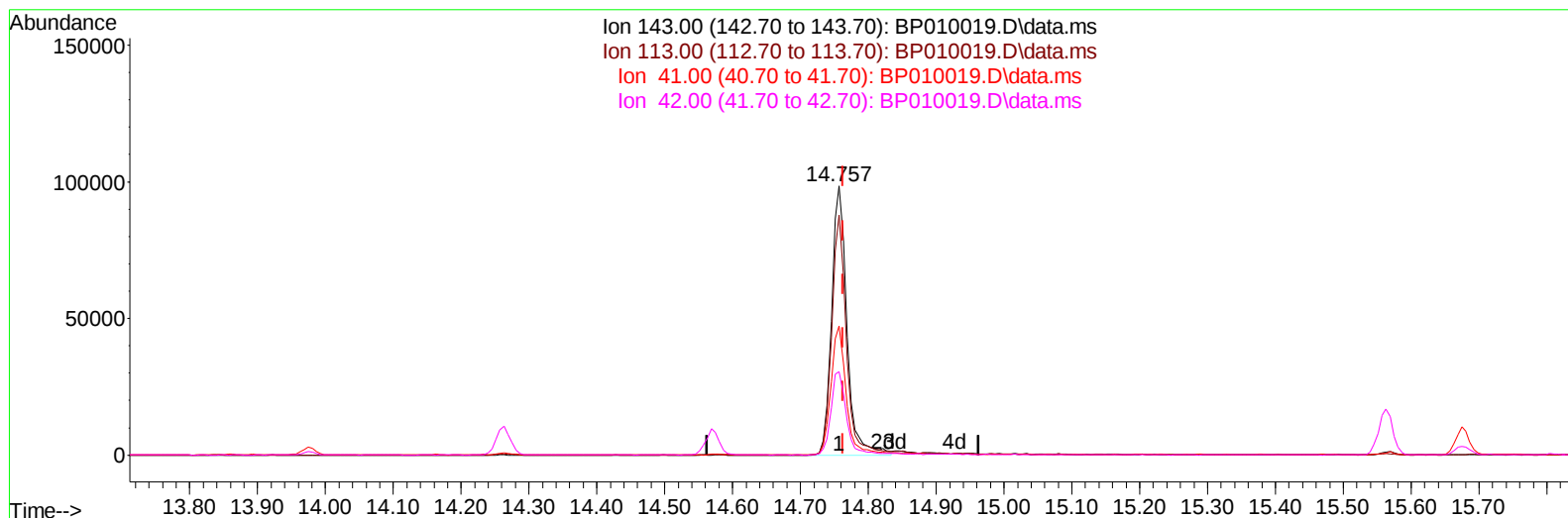
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 Data File : BP010019.D
 Acq On : 21 Apr 2022 17:53
 Operator : CG/JU
 Sample : PB144260BL
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK260

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TIC: BP010019.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.757min (-0.006) 27.12 ng/ul m

response 153225

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	132.70	89.21#
41.00	23.20	48.07#
42.00	18.00	30.93#

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 Sample : PB144260BL
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 ALS Vial : 14 Sample Multiplier: 1

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.940	152	117138	20.000	ng/ul	0.00
20) Naphthalene-d8	10.746	136	525254	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.575	164	394525	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.322	188	885640	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.404	240	841023	20.000	ng/ul	# 0.00
88) Perylene-d12	23.857	264	742023	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.369	96	17650	6.571	ng/uL	0.00
4) Pyridine-d5	3.787	84	232301	30.740	ng/ul	0.00
7) Phenol-d5	7.099	99	297896	27.959	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.263	67	209059	32.700	ng/ul	0.00
11) 2-Chlorophenol-d4	7.469	132	241605	30.618	ng/ul	0.00
15) 4-Methylphenol-d8	8.646	113	265172	30.038	ng/ul	0.00
21) Nitrobenzene-d5	9.099	128	132453	31.706	ng/ul	0.00
24) 2-Nitrophenol-d4	9.822	143	147107	32.243	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.363	165	255052m	28.716	ng/ul	0.00
31) 4-Chloroaniline-d4	10.875	131	377099	29.610	ng/ul	0.00
46) Dimethylphthalate-d6	13.975	166	1039853	33.883	ng/ul	0.00
49) Acenaphthylene-d8	14.263	160	1119014	30.127	ng/ul	0.00
54) 4-Nitrophenol-d4	14.757	143	153225m	27.120	ng/ul	0.00
60) Fluorene-d10	15.563	176	880503	33.353	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.675	200	144609	25.823	ng/ul	0.00
73) Anthracene-d10	17.422	188	1406587	32.829	ng/ul	0.00
81) Pyrene-d10	19.651	212	1698500	35.512	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.698	264	1303234	33.501	ng/ul	0.00

Target Compounds Qvalue

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