

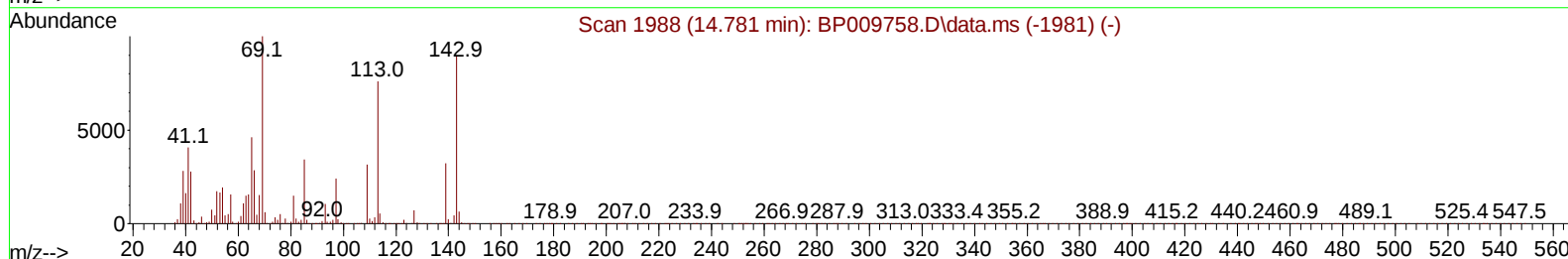
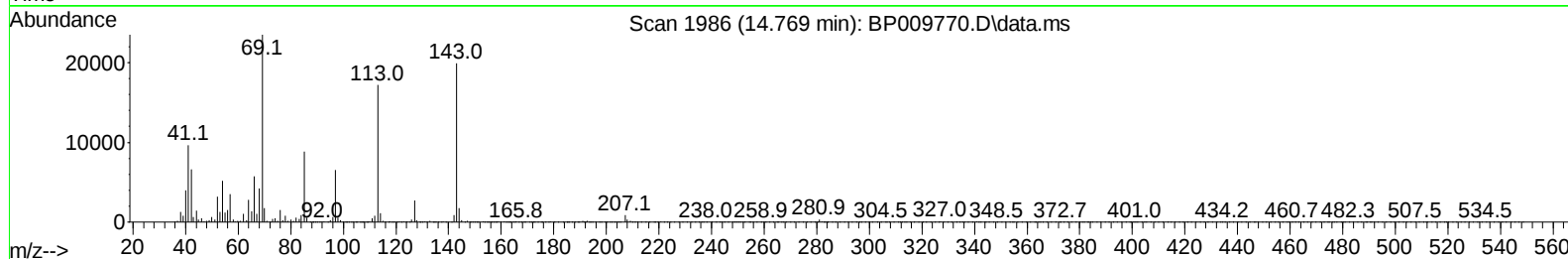
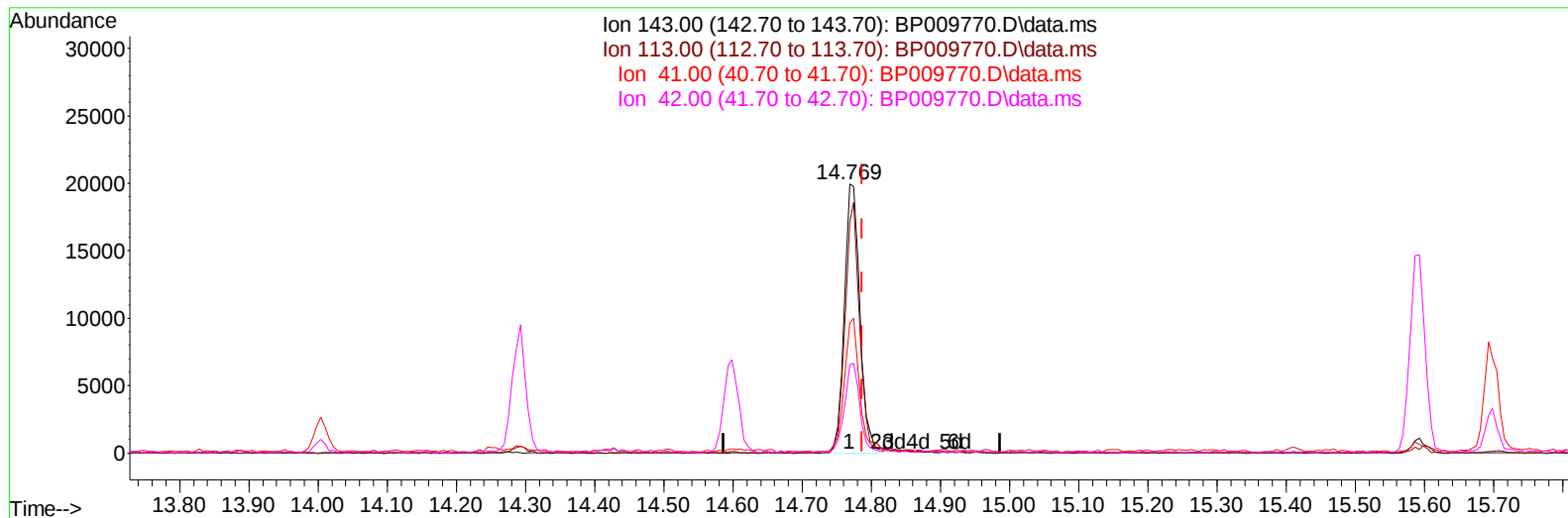
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041122\
 Data File : BP009770.D
 Acq On : 12 Apr 2022 02:04
 Operator : CG/JU
 Sample : N2338-16
 Misc :
 ALS Vial : 27 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0J88

Manual IntegrationsAPPROVED

Quant Time: Apr 12 02:45:50 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP040722.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Apr 07 19:16:52 2022
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 04/13/2022
 Supervised By :Yogesh Patel 04/13/2022



TIC: BP009770.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.769min (-0.018) 6.92 ng/u1

response 31052

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	132.70	86.24#
41.00	23.20	48.55#
42.00	18.00	33.02#

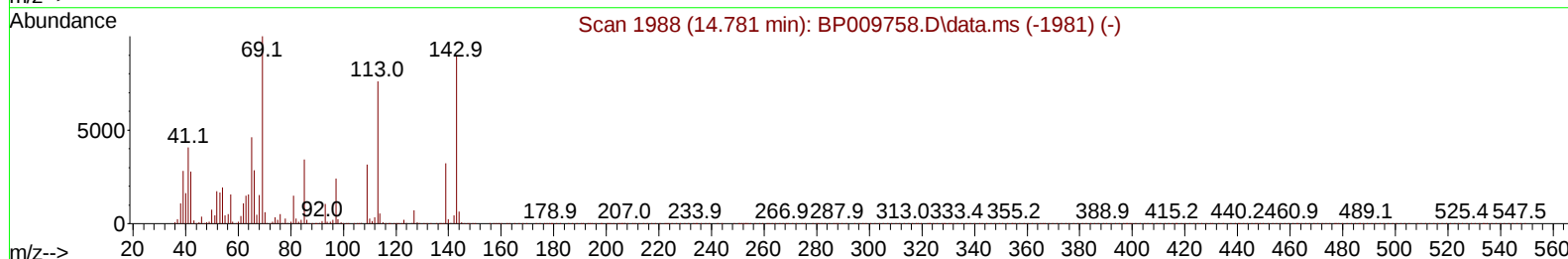
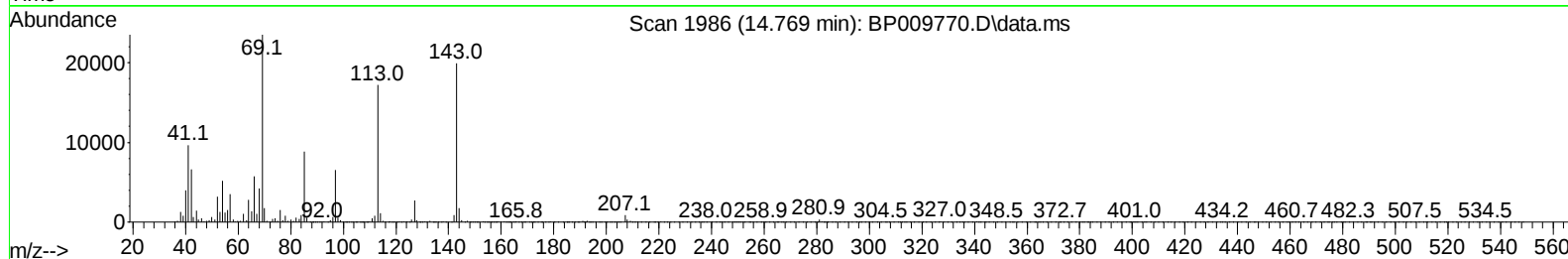
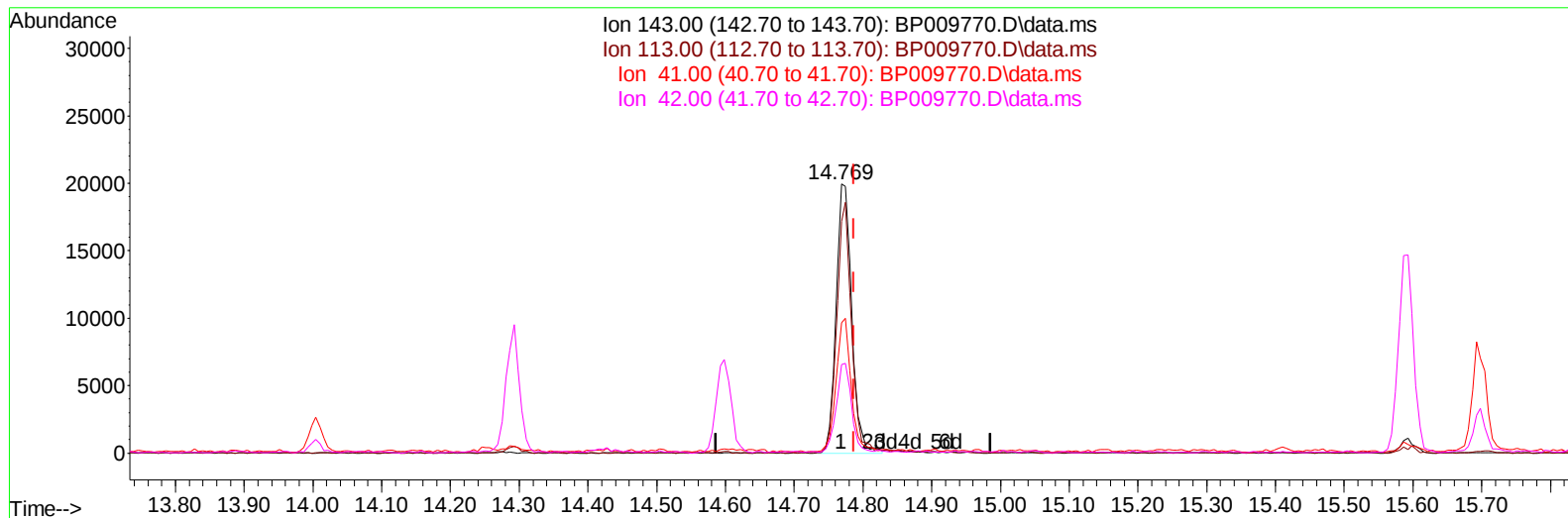
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TIC: BP009770.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.769min (-0.018) 7.00 ng/ul m

response 31393

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	132.70	86.24#
41.00	23.20	48.55#
42.00	18.00	33.02#

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.969	152	95415	20.000	ng/ul	-0.01
20) Naphthalene-d8	10.775	136	432366	20.000	ng/ul	-0.02
38) Acenaphthene-d10	14.598	164	314961	20.000	ng/ul	-0.01
64) Phenanthrene-d10	17.351	188	713696	20.000	ng/ul	#-0.01
79) Chrysene-d12	21.427	240	816281	20.000	ng/ul	#-0.02
88) Perylene-d12	23.898	264	850840	20.000	ng/ul	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.387	96	12963	6.063	ng/uL	0.00
4) Pyridine-d5	3.811	84	34998	5.937	ng/ul	0.00
7) Phenol-d5	7.122	99	55187	6.588	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.293	67	171643	34.418	ng/ul	-0.01
11) 2-Chlorophenol-d4	7.499	132	154843	24.505	ng/ul	0.00
15) 4-Methylphenol-d8	8.669	113	109237	15.621	ng/ul	-0.01
21) Nitrobenzene-d5	9.128	128	111705	35.827	ng/ul	-0.01
24) 2-Nitrophenol-d4	9.851	143	106695	31.213	ng/ul	-0.01
28) 2,4-Dichlorophenol-d3	10.393	165	186889	26.127	ng/ul	-0.01
31) 4-Chloroaniline-d4	10.904	131	204243	20.047	ng/ul	-0.01
46) Dimethylphthalate-d6	14.004	166	874528	35.348	ng/ul	-0.02
49) Acenaphthylene-d8	14.292	160	939359	32.048	ng/ul	-0.01
54) 4-Nitrophenol-d4	14.769	143	31393m	6.997	ng/ul	-0.02
60) Fluorene-d10	15.592	176	762975	36.659	ng/ul	-0.01
65) 4,6-Dinitro-2-methylph...	15.698	200	122096	28.878	ng/ul	-0.01
73) Anthracene-d10	17.451	188	1382350	40.454	ng/ul	-0.01
81) Pyrene-d10	19.674	212	1720439	38.591	ng/ul	-0.01
92) Benzo(a)pyrene-d12	23.739	264	1825459	41.199	ng/ul	-0.02

Target Compounds Qvalue

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

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