

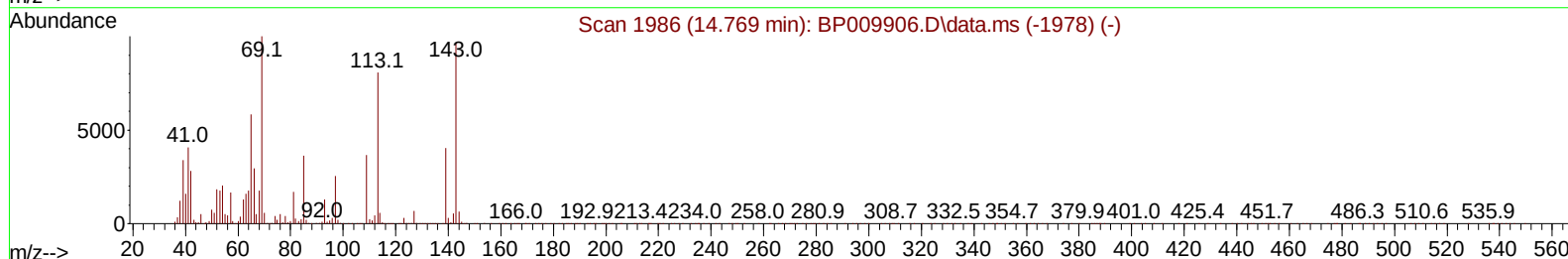
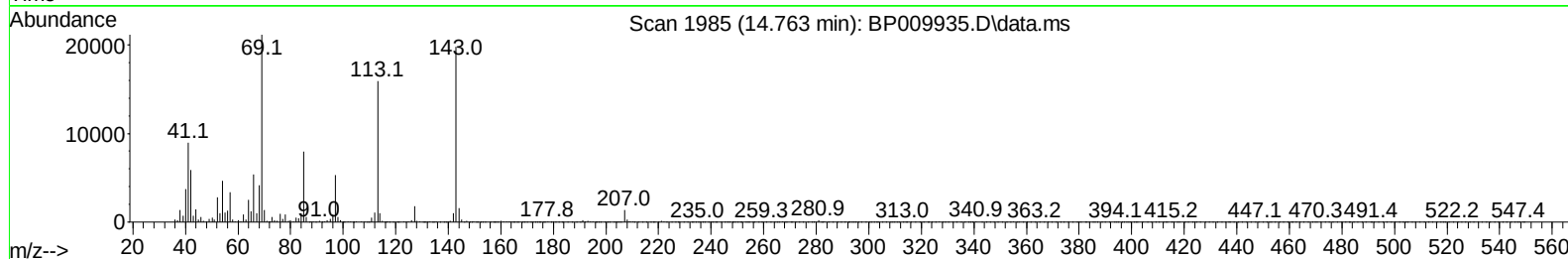
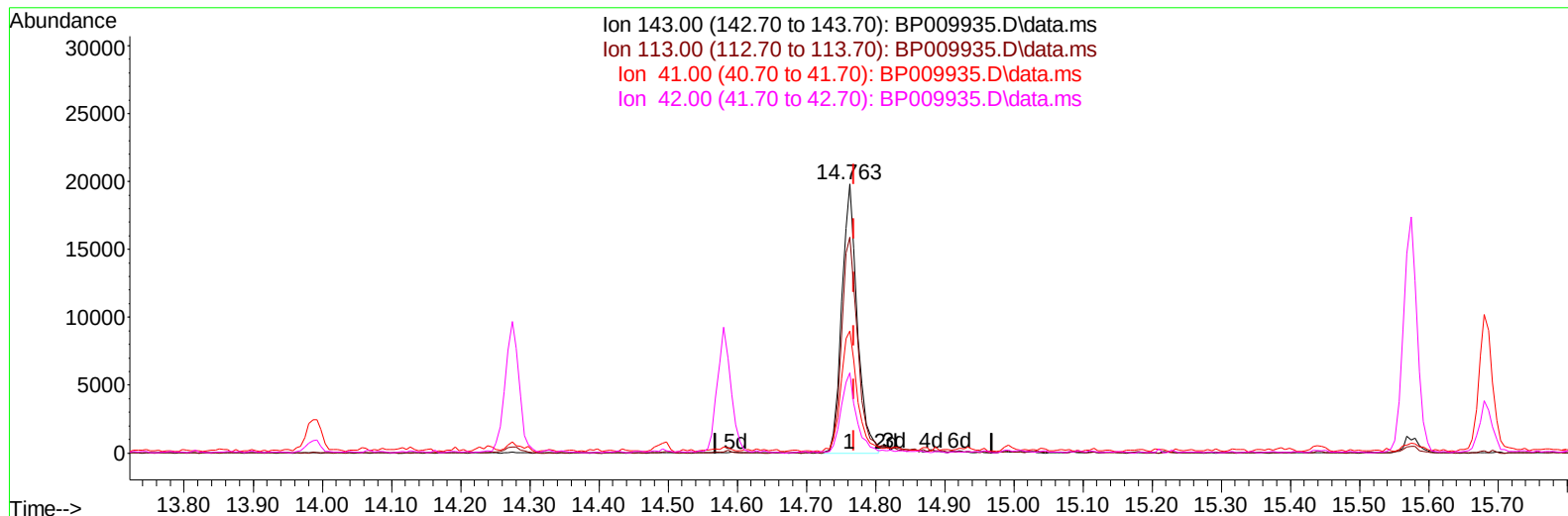
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041822\
 Data File : BP009935.D
 Acq On : 19 Apr 2022 02:14
 Operator : CG/JU
 Sample : N2421-18
 Misc :
 ALS Vial : 30 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0J80

Manual IntegrationsAPPROVED

Quant Time: Apr 19 02:44:38 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP040722.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Apr 18 00:59:30 2022
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 04/19/2022
 Supervised By :Yogesh Patel 04/22/2022



TIC: BP009935.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.763min (-0.006) 5.29 ng/ul

response 30868

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	132.70	80.24#
41.00	23.20	45.43#
42.00	18.00	29.74#

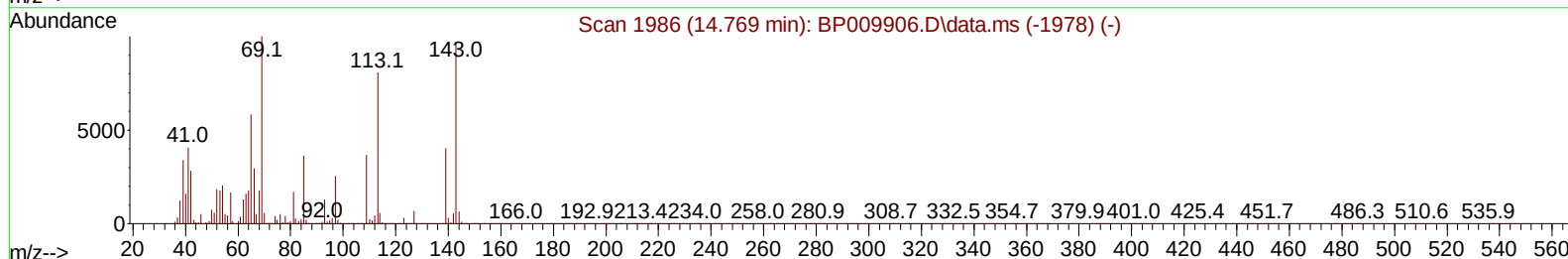
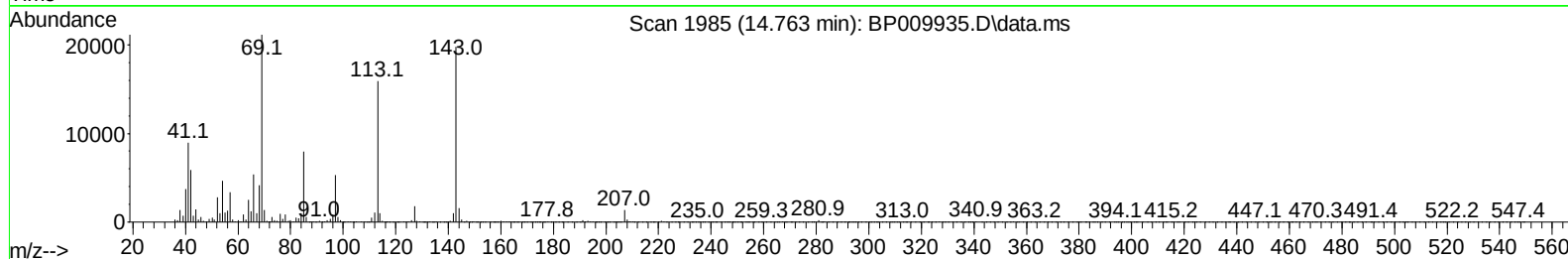
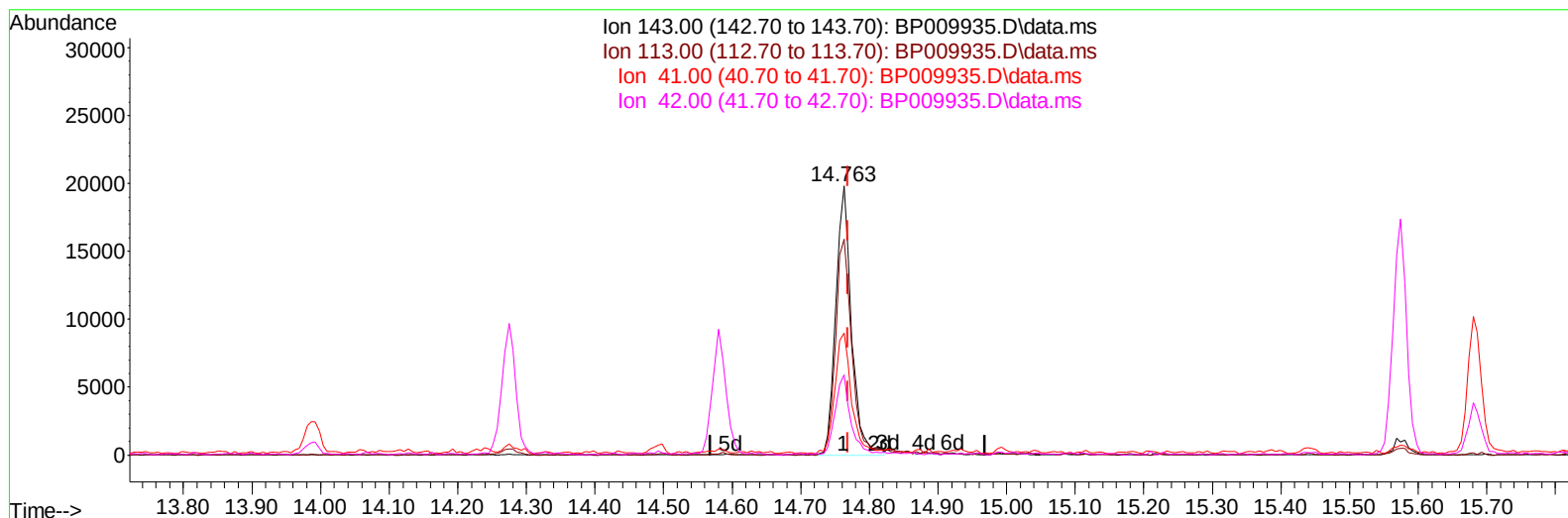
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TIC: BP009935.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.763min (-0.006) 5.37 ng/ul m

response 31336

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	132.70	80.24#
41.00	23.20	45.43#
42.00	18.00	29.74#

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.951	152	125117	20.000	ng/ul	0.00
20) Naphthalene-d8	10.757	136	565853	20.000	ng/ul	-0.01
38) Acenaphthene-d10	14.580	164	409602	20.000	ng/ul	-0.01
64) Phenanthrene-d10	17.333	188	904928	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.415	240	905327	20.000	ng/ul	# 0.00
88) Perylene-d12	23.880	264	818668	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.381	96	12079	4.309	ng/uL	0.00
4) Pyridine-d5	3.799	84	68441	8.855	ng/ul	0.00
7) Phenol-d5	7.104	99	70934	6.458	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.275	67	202915	31.029	ng/ul	0.00
11) 2-Chlorophenol-d4	7.481	132	191446	23.105	ng/ul	0.00
15) 4-Methylphenol-d8	8.651	113	143382	15.636	ng/ul	-0.01
21) Nitrobenzene-d5	9.110	128	133756	32.779	ng/ul	0.00
24) 2-Nitrophenol-d4	9.834	143	134031	29.960	ng/ul	-0.01
28) 2,4-Dichlorophenol-d3	10.375	165	236057	25.216	ng/ul	0.00
31) 4-Chloroaniline-d4	10.886	131	387679	29.076	ng/ul	0.00
46) Dimethylphthalate-d6	13.986	166	1014567	31.533	ng/ul	-0.01
49) Acenaphthylene-d8	14.275	160	1103575	28.951	ng/ul	-0.01
54) 4-Nitrophenol-d4	14.763	143	31336m	5.370	ng/ul	0.00
60) Fluorene-d10	15.574	176	887195	32.778	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.686	200	150974	28.162	ng/ul	0.00
73) Anthracene-d10	17.433	188	1546151	35.685	ng/ul	0.00
81) Pyrene-d10	19.663	212	1904178	38.511	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.721	264	1602187	37.581	ng/ul	0.00

Target Compounds Qvalue

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

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