

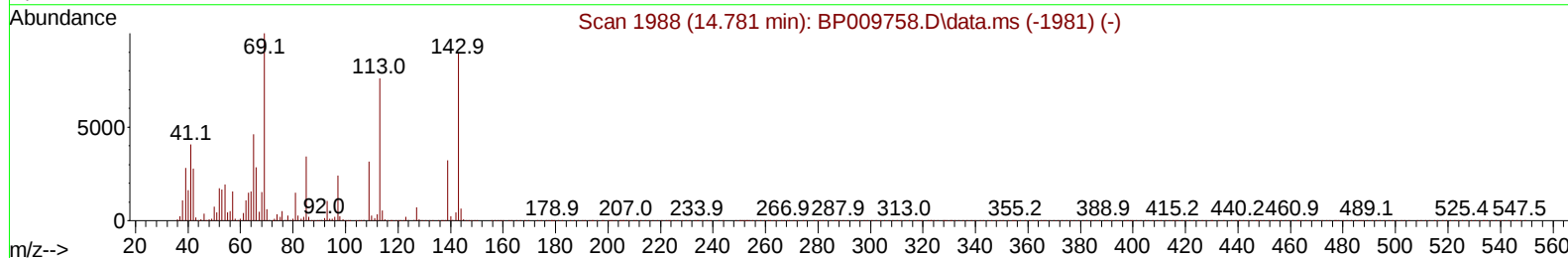
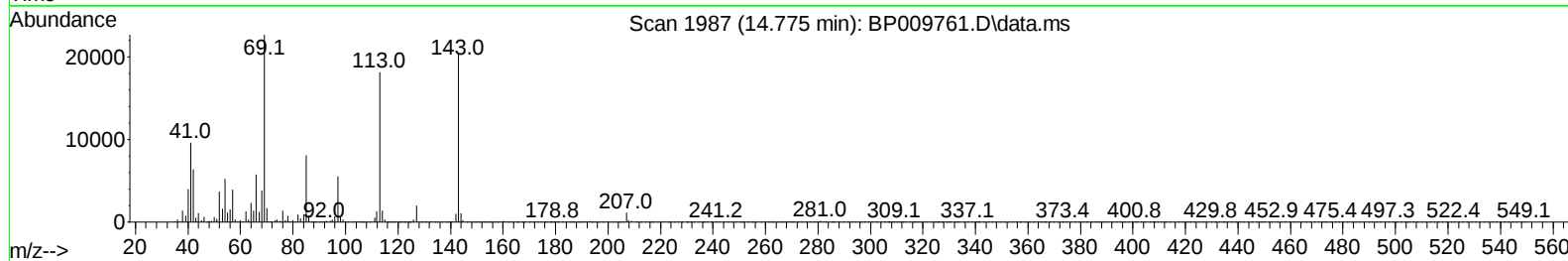
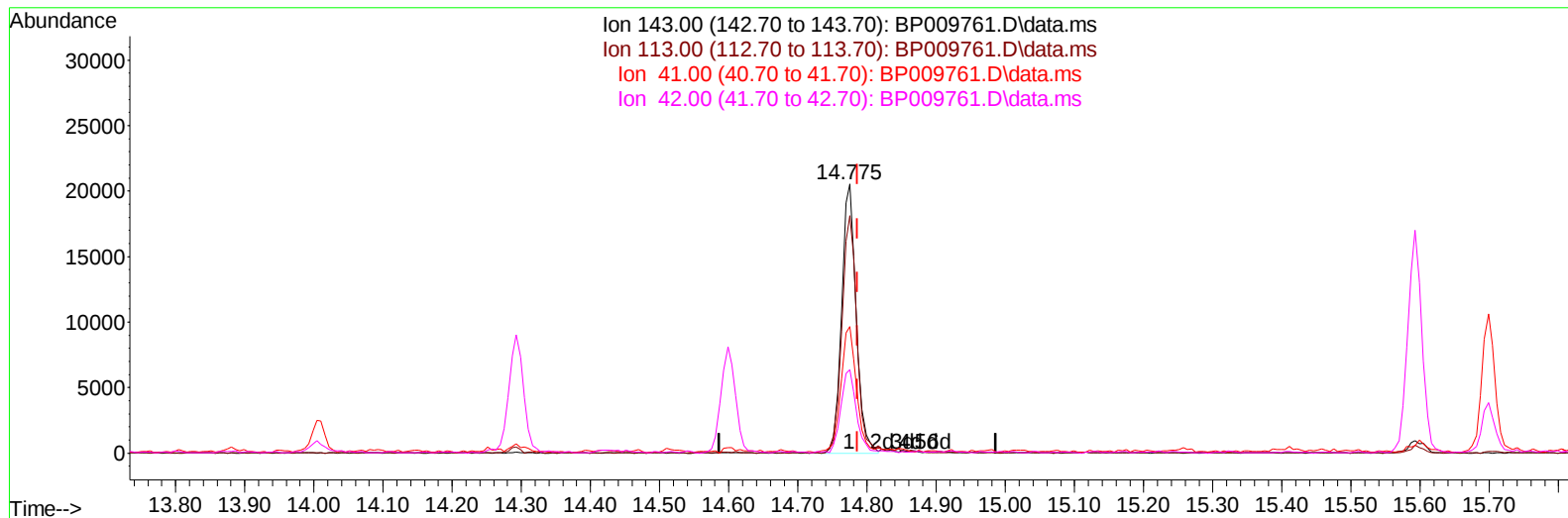
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041122\
 Data File : BP009761.D
 Acq On : 11 Apr 2022 21:01
 Operator : CG/JU
 Sample : N2338-14
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0J79

Manual IntegrationsAPPROVED

Quant Time: Apr 12 01:53:23 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: BP009761.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.775min (-0.011) 6.64 ng/u1

response 29964

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	132.70	88.39#
41.00	23.20	47.04#
42.00	18.00	31.05#

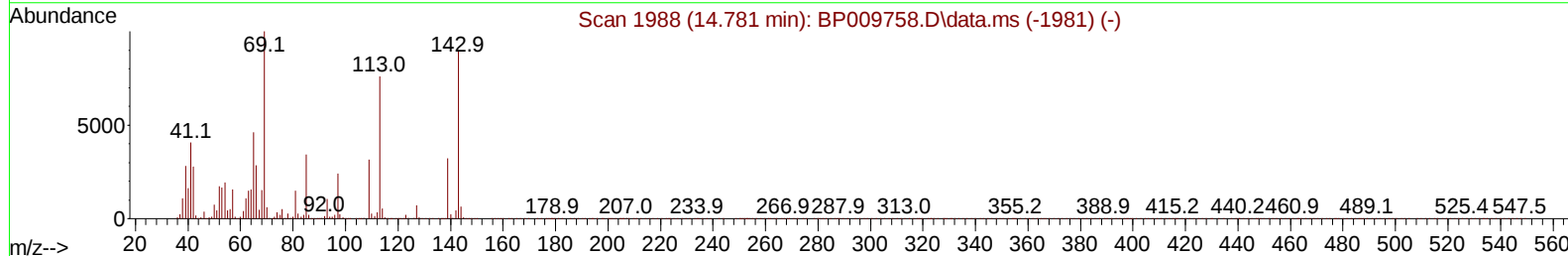
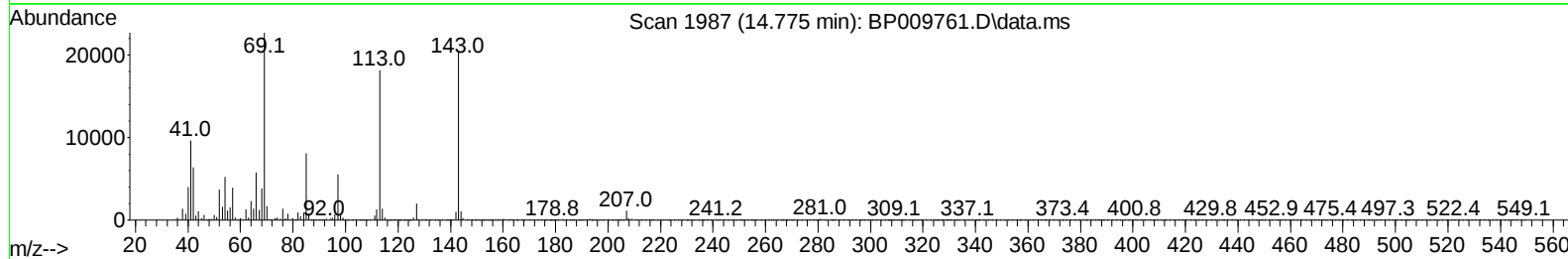
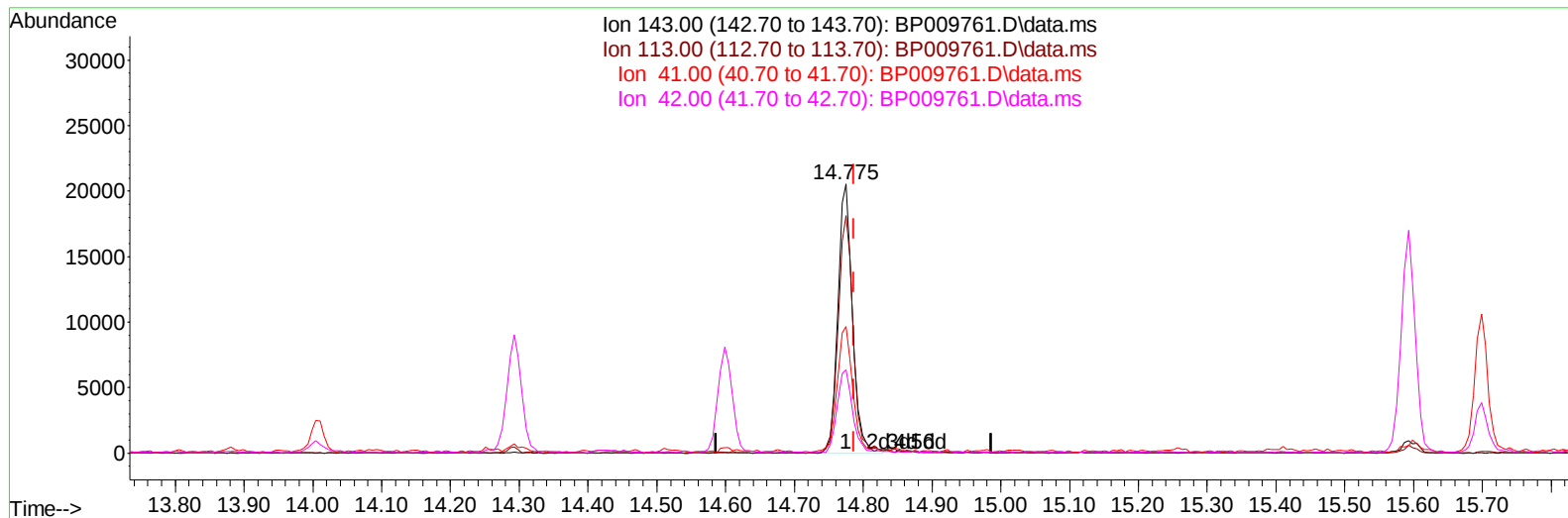
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TIC: BP009761.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.775min (-0.011) 6.68 ng/ul m

response 30127

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	132.70	88.39#
41.00	23.20	47.04#
42.00	18.00	31.05#

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 Operator : CG/JU
 Sample : N2338-14
 Misc :
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Instrument :
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.969	152	102620	20.000	ng/ul	-0.01
20) Naphthalene-d8	10.775	136	440409	20.000	ng/ul	-0.02
38) Acenaphthene-d10	14.599	164	316674	20.000	ng/ul	-0.01
64) Phenanthrene-d10	17.351	188	706180	20.000	ng/ul	#-0.01
79) Chrysene-d12	21.433	240	780899	20.000	ng/ul	#-0.01
88) Perylene-d12	23.904	264	830718	20.000	ng/ul	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.387	96	14384	6.256	ng/uL	0.00
4) Pyridine-d5	3.811	84	16566	2.613	ng/ul	0.00
7) Phenol-d5	7.122	99	53360	5.923	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.293	67	176832	32.969	ng/ul	-0.01
11) 2-Chlorophenol-d4	7.499	132	165620	24.371	ng/ul	0.00
15) 4-Methylphenol-d8	8.669	113	108950	14.486	ng/ul	-0.01
21) Nitrobenzene-d5	9.128	128	112231	35.338	ng/ul	-0.01
24) 2-Nitrophenol-d4	9.852	143	113723	32.661	ng/ul	-0.01
28) 2,4-Dichlorophenol-d3	10.393	165	194099	26.640	ng/ul	-0.01
31) 4-Chloroaniline-d4	10.905	131	196682	18.953	ng/ul	-0.01
46) Dimethylphthalate-d6	14.010	166	862728	34.682	ng/ul	-0.01
49) Acenaphthylene-d8	14.293	160	935805	31.754	ng/ul	-0.01
54) 4-Nitrophenol-d4	14.775	143	30127m	6.678	ng/ul	-0.01
60) Fluorene-d10	15.593	176	766693	36.638	ng/ul	-0.01
65) 4,6-Dinitro-2-methylph...	15.698	200	148506	35.499	ng/ul	-0.01
73) Anthracene-d10	17.451	188	1406847	41.609	ng/ul	-0.01
81) Pyrene-d10	19.681	212	1758081	41.222	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.745	264	1869846	43.223	ng/ul	-0.02

Target Compounds Qvalue

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

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