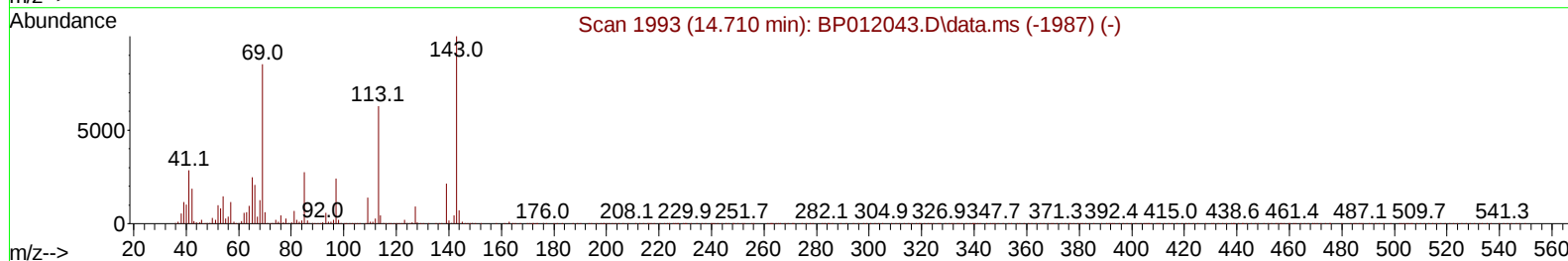
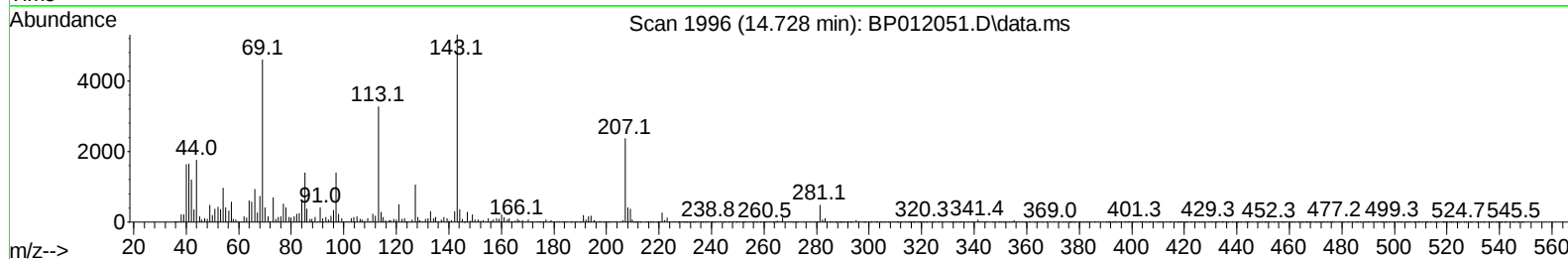
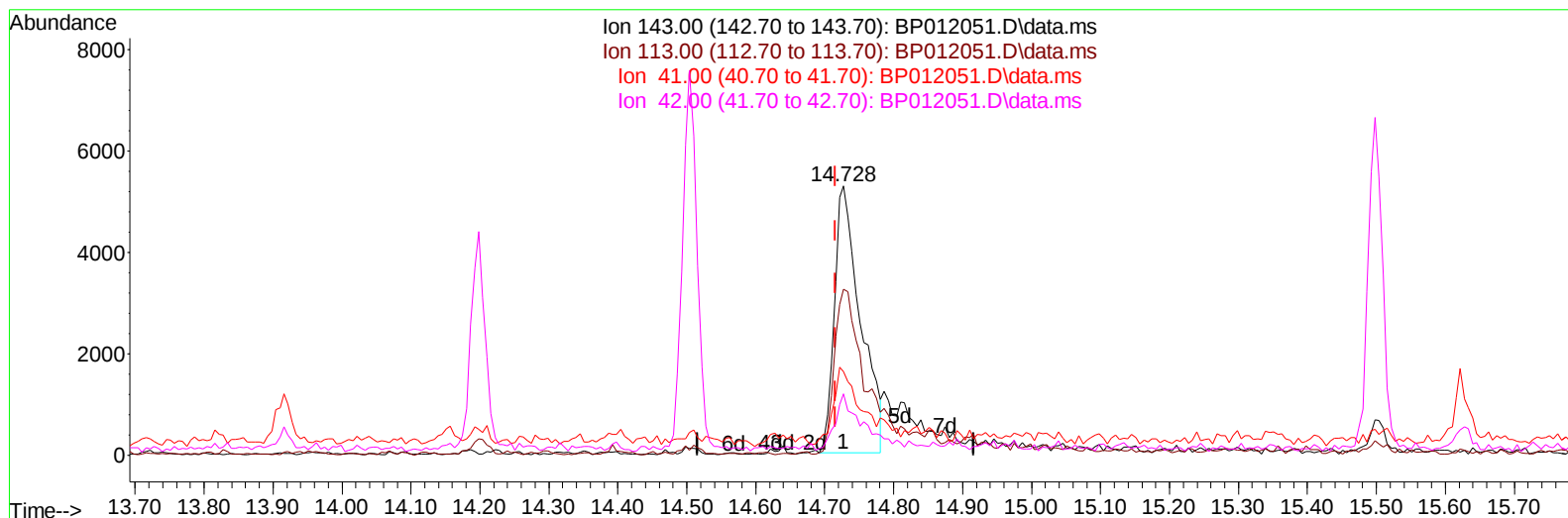


Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101122\
Data File : BP012051.D
Acq On : 11 Oct 2022 21:25
Operator : CG/JU
Sample : N4991-01
Misc :
ALS Vial : 19 Sample Multiplier: 1

Quant Time: Oct 12 00:24:40 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP101122.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue Oct 11 23:54:43 2022
Response via : Initial Calibration



TIC: BP012051.D\data.ms

(54) 4-Nitrophenol-d4 (S)

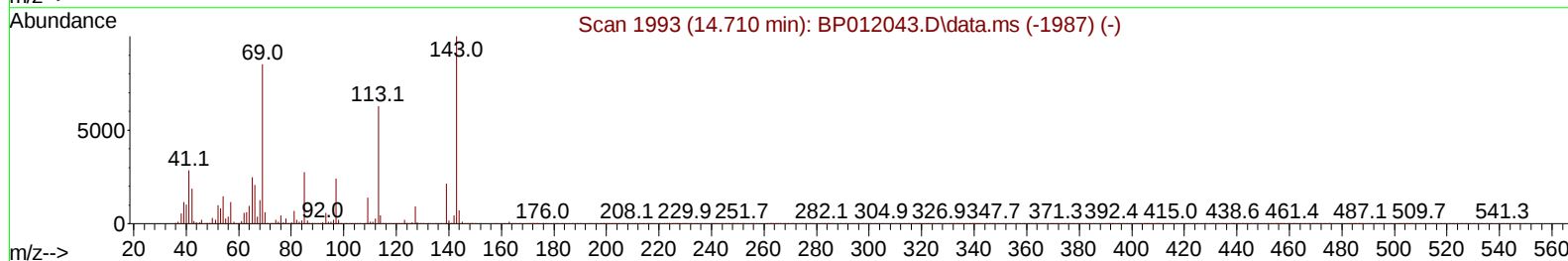
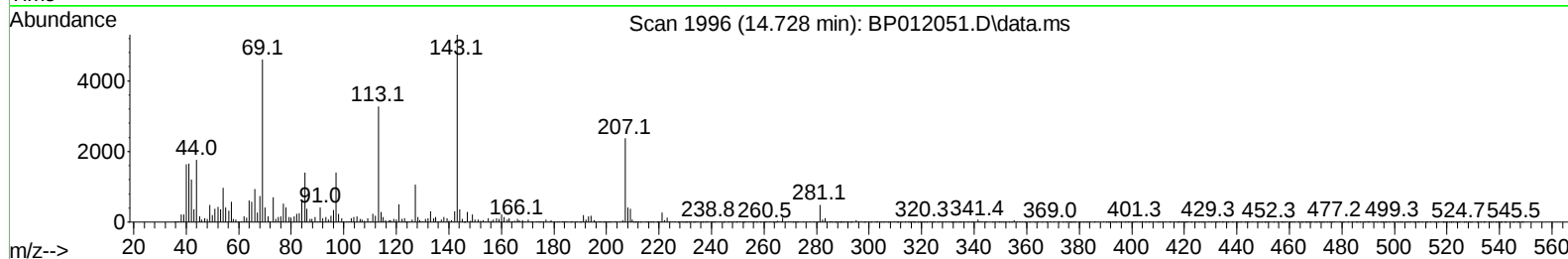
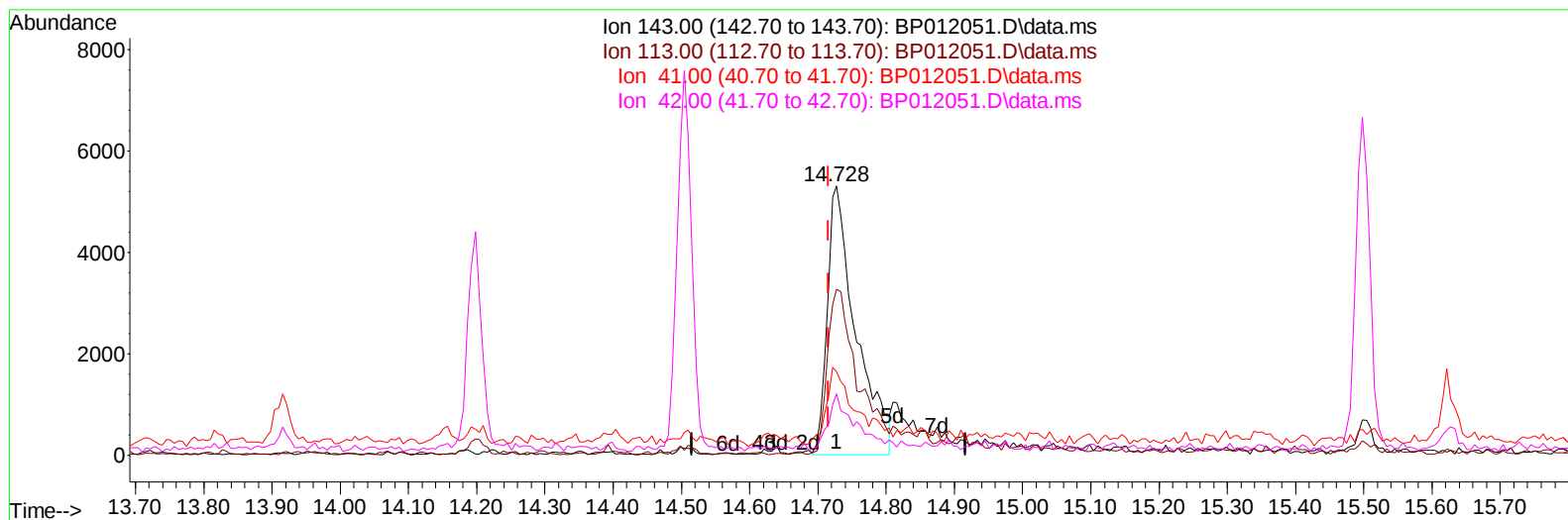
14.728min (+ 0.012) 1.91 ng/ul

response 13800

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	60.80	61.76
41.00	26.60	31.12
42.00	17.60	22.88#

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101122\
Data File : BP012051.D
Acq On : 11 Oct 2022 21:25
Operator : CG/JU
Sample : N4991-01
Misc :
ALS Vial : 19 Sample Multiplier: 1

Quant Time: Oct 12 00:24:40 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP101122.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue Oct 11 23:54:43 2022
Response via : Initial Calibration



TIC: BP012051.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.728min (+ 0.012) 2.13 ng/ul m

response 15379

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	60.80	61.76
41.00	26.60	31.12
42.00	17.60	22.88#

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101122\
 Data File : BP012051.D
 Acq On : 11 Oct 2022 21:25
 Operator : CG/JU
 Sample : N4991-01
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Quant Time: Oct 12 00:24:40 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP101122.M
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 QLast Update : Tue Oct 11 23:54:43 2022
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.857	152	267169	20.000	ng/ul	0.00
20) Naphthalene-d8	10.663	136	1073376	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.504	164	579927	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.263	188	1076540	20.000	ng/ul	0.00
79) Chrysene-d12	21.351	240	1081403	20.000	ng/ul	0.00
88) Perylene-d12	23.768	264	1288638	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.281	96	20478	3.022	ng/uL	0.00
4) Pyridine-d5	0.000	84	0d	0.000	ng/ul	
7) Phenol-d5	7.022	99	253090	11.779	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.193	67	206706	16.948	ng/ul	0.00
11) 2-Chlorophenol-d4	7.387	132	263197	15.431	ng/ul	0.00
15) 4-Methylphenol-d8	8.569	113	164761	9.983	ng/ul	0.00
21) Nitrobenzene-d5	9.028	128	134552	17.749	ng/ul	0.00
24) 2-Nitrophenol-d4	9.751	143	127951	16.646	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.287	165	206391	13.854	ng/ul	0.00
31) 4-Chloroaniline-d4	10.810	131	221380	10.084	ng/ul	0.00
46) Dimethylphthalate-d6	13.916	166	639323	17.407	ng/ul	0.00
49) Acenaphthylene-d8	14.198	160	778782	17.148	ng/ul	0.00
54) 4-Nitrophenol-d4	14.728	143	15379m	2.133	ng/ul	0.01
60) Fluorene-d10	15.498	176	589961	17.754	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.622	200	25977	4.570	ng/ul	0.00
73) Anthracene-d10	17.363	188	833163	17.823	ng/ul	0.00
81) Pyrene-d10	19.598	212	950617	17.159	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.609	264	879058	14.033	ng/ul	0.00

Target Compounds	Qvalue

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

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