

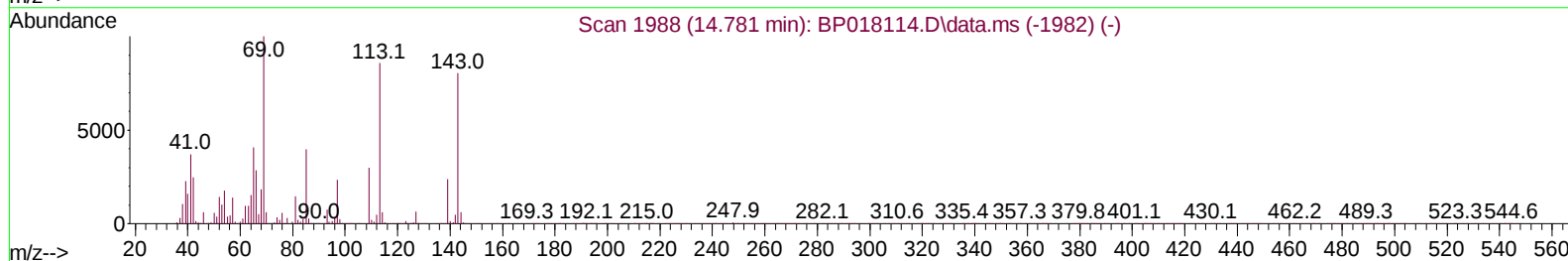
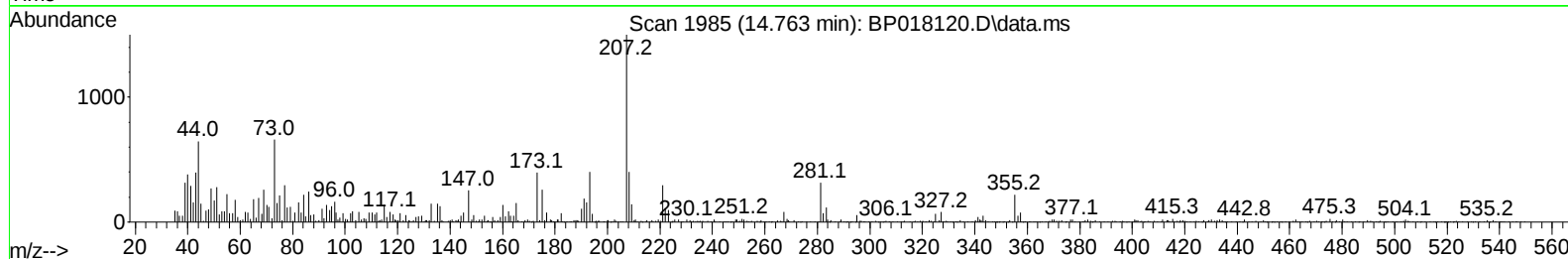
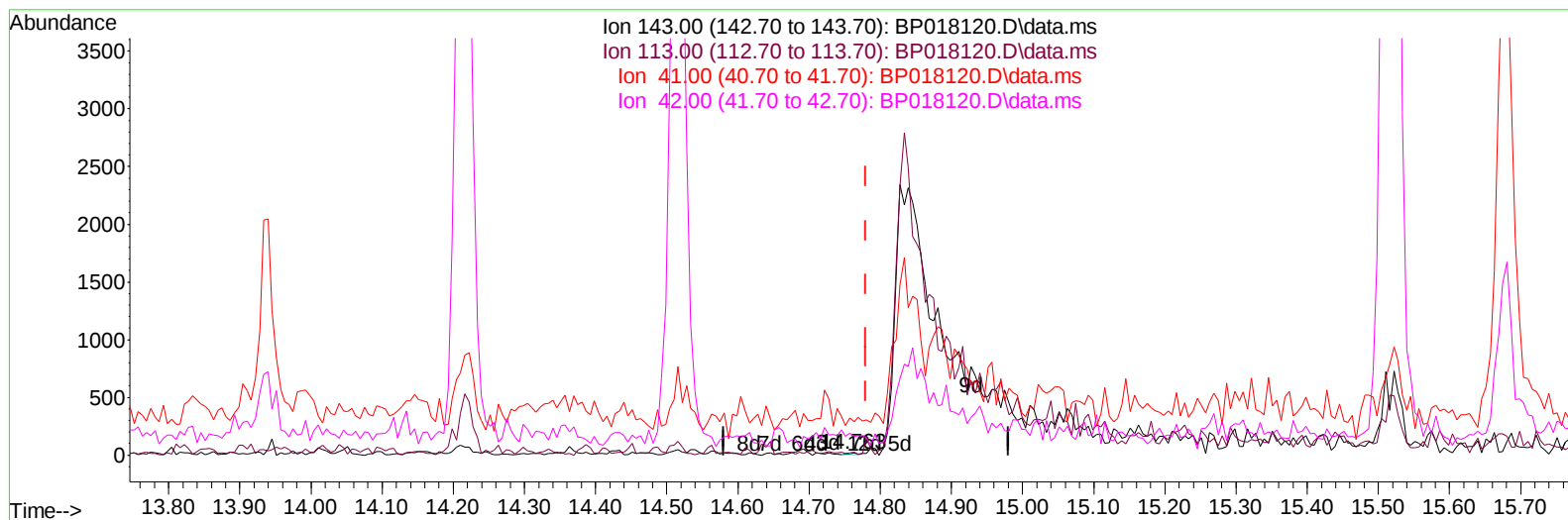
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111023\
 Data File : BP018120.D
 Acq On : 13 Nov 2023 00:31
 Operator : MA/JU
 Sample : 08082-06
 Misc :
 ALS Vial : 101 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BH332

Manual IntegrationsAPPROVED

Quant Time: Nov 13 03:25:31 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP111023.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Nov 10 21:55:18 2023
 Response via : Initial Calibration

Reviewed By :Yogesh Patel 11/14/2023
 Supervised By :mohammad ahmed 11/15/2023



TIC: BP018120.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.763min (-0.017) 0.00 ng/u1

response 10

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	103.00	131.58#
41.00	47.70	1526.32#
42.00	30.10	842.11#

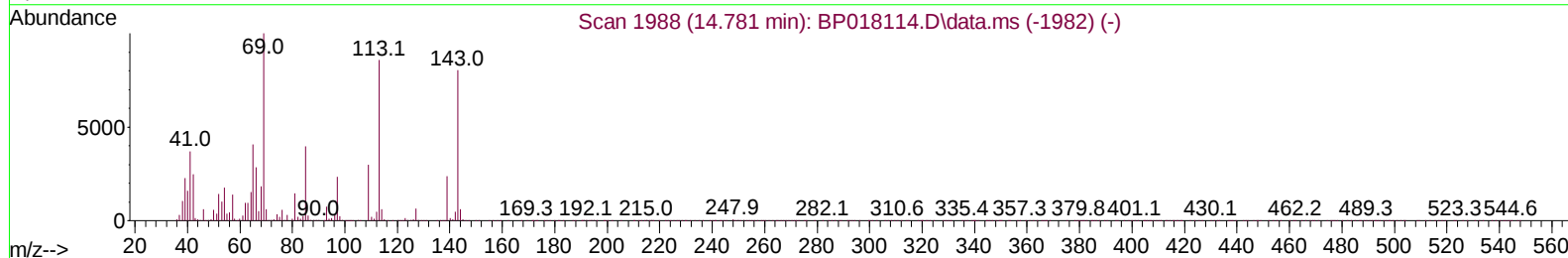
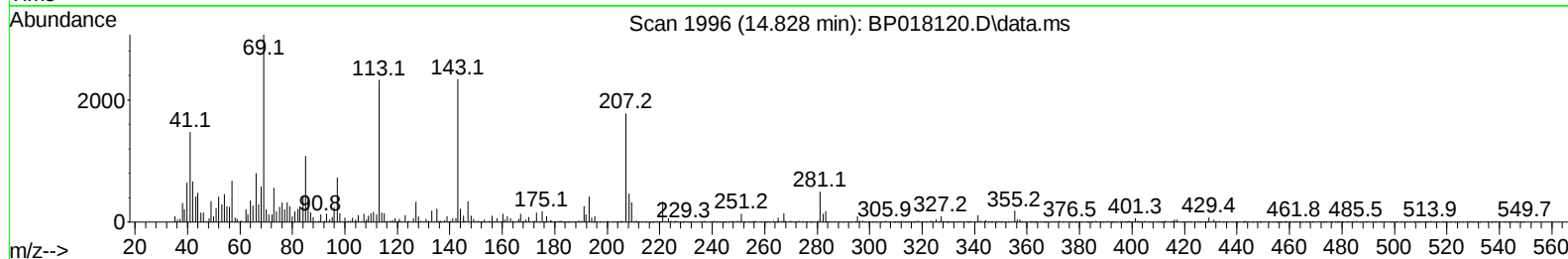
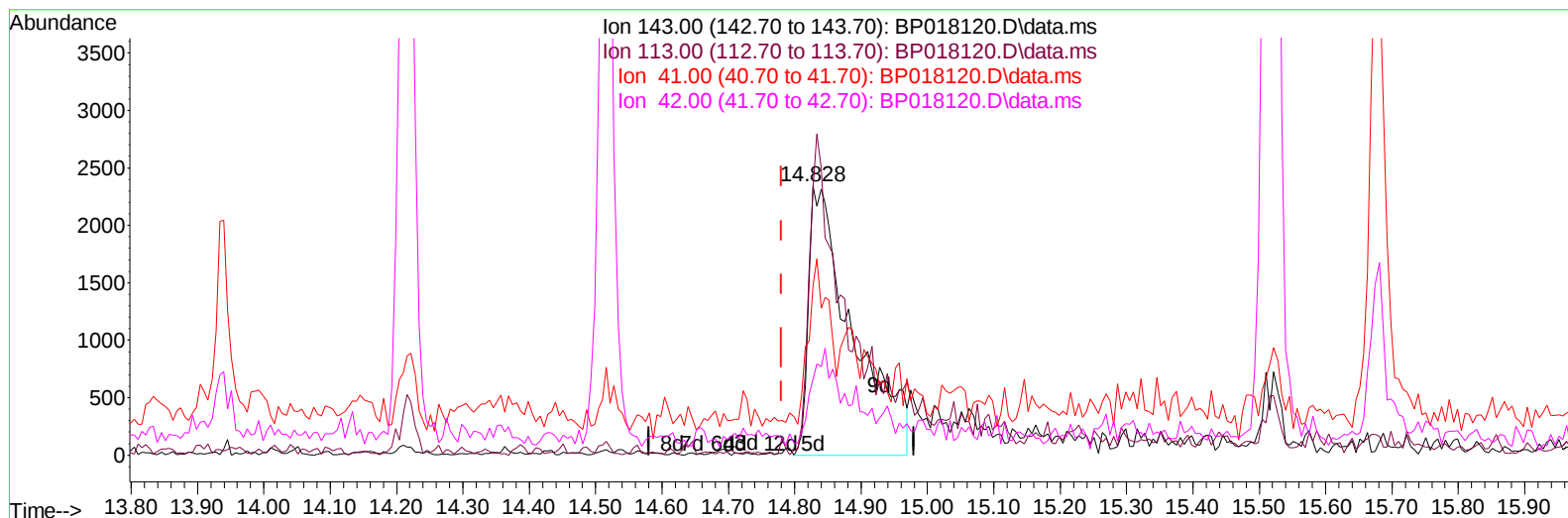
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TIC: BP018120.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.828min (+ 0.047) 3.93 ng/ul m

response 10785

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	103.00	99.53
41.00	47.70	63.09#
42.00	30.10	28.28

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Quant Time: Nov 13 03:26:47 2023
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.840	152	79771	20.000	ng/ul	-0.01
20) Naphthalene-d8	10.663	136	325562	20.000	ng/ul	-0.01
38) Acenaphthene-d10	14.516	164	199218	20.000	ng/ul	-0.01
64) Phenanthrene-d10	17.298	188	379097	20.000	ng/ul	-0.01
79) Chrysene-d12	21.422	240	325048	20.000	ng/ul	-0.01
88) Perylene-d12	23.927	264	381202	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.228	96	8520	3.768	ng/uL	0.00
4) Pyridine-d5	3.664	84	63209	10.720	ng/ul	0.00
7) Phenol-d5	7.022	99	60160	7.812	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.193	67	135777	30.000	ng/ul	0.00
11) 2-Chlorophenol-d4	7.369	132	151361	24.819	ng/ul	0.00
15) 4-Methylphenol-d8	8.575	113	106901	16.940	ng/ul	0.00
21) Nitrobenzene-d5	9.046	128	111450	37.997	ng/ul	-0.01
24) 2-Nitrophenol-d4	9.763	143	98095	29.520	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.299	165	149521	25.246	ng/ul	0.00
31) 4-Chloroaniline-d4	10.840	131	198677	24.462	ng/ul	-0.01
46) Dimethylphthalate-d6	13.940	166	565923	30.788	ng/ul	-0.01
49) Acenaphthylene-d8	14.216	160	607025	30.320	ng/ul	-0.01
54) 4-Nitrophenol-d4	14.828	143	10785m	3.926	ng/ul	0.05
60) Fluorene-d10	15.522	176	460276	30.329	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.681	200	66811	24.378	ng/ul	0.00
73) Anthracene-d10	17.398	188	666428	32.504	ng/ul	-0.01
81) Pyrene-d10	19.651	212	744689	34.299	ng/ul	-0.01
92) Benzo(a)pyrene-d12	23.757	264	763573	34.448	ng/ul	-0.02

Target Compounds Qvalue

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

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