

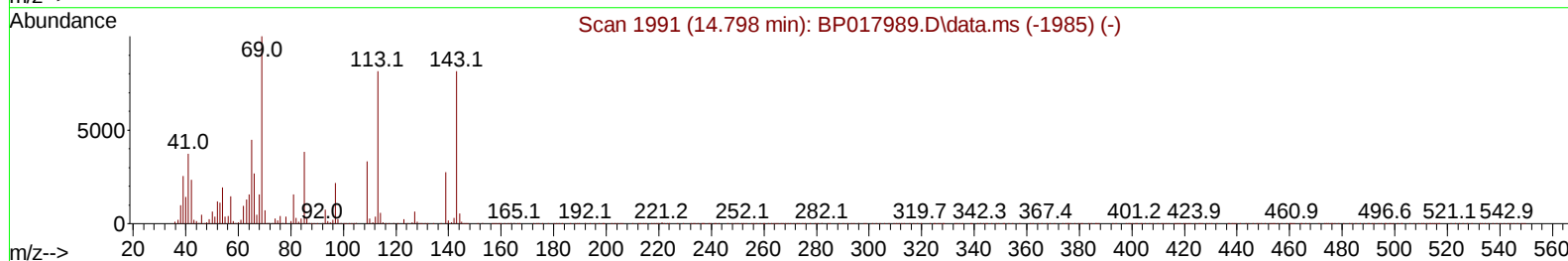
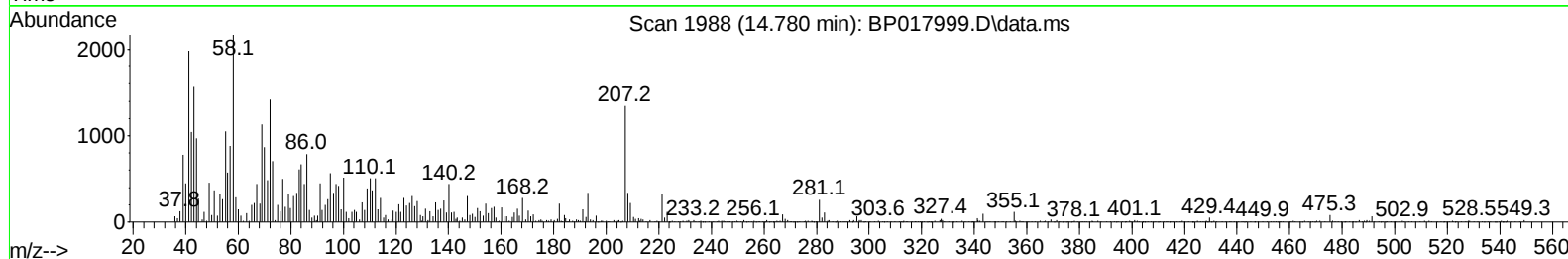
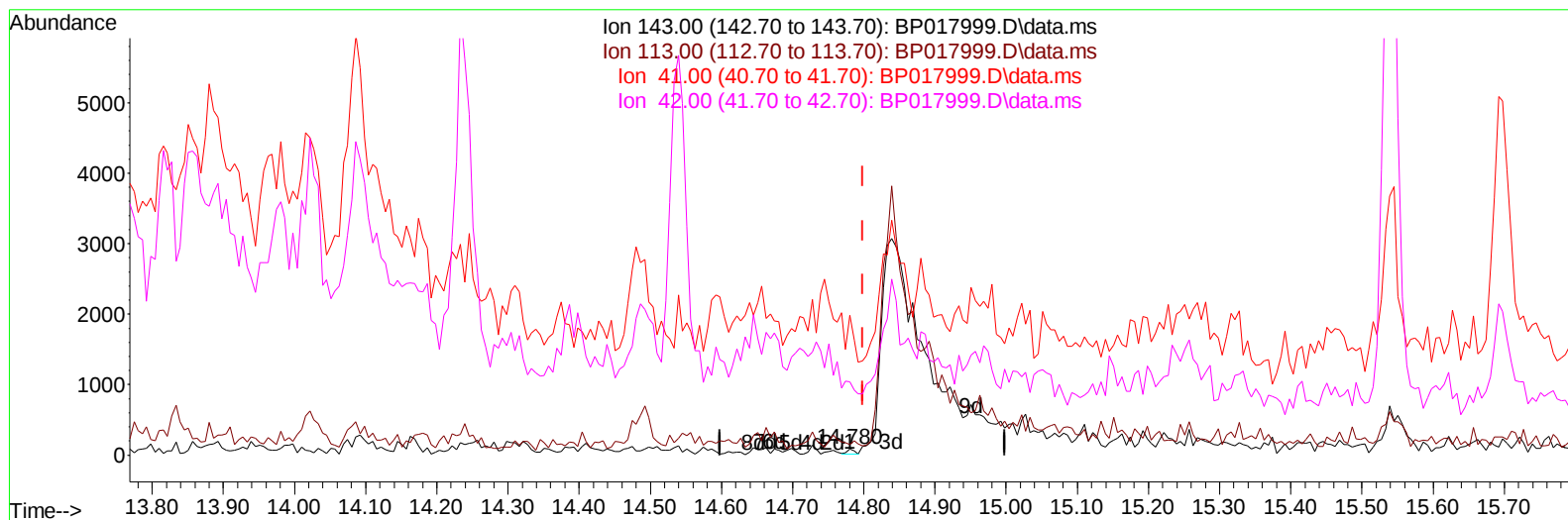
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110923\
Data File : BP017999.D
Acq On : 09 Nov 2023 21:15
Operator : MA/JU
Sample : 05082-07
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BH358

Manual IntegrationsAPPROVED

Quant Time: Nov 09 23:18:27 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP110923.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Thu Nov 09 22:34:38 2023
Response via : Initial Calibration

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



TIC: BP017999.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.780min (-0.018) 0.02 ng/u1

response 46

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	100.40	167.42#
41.00	46.90	2230.34#
42.00	29.70	1171.91#

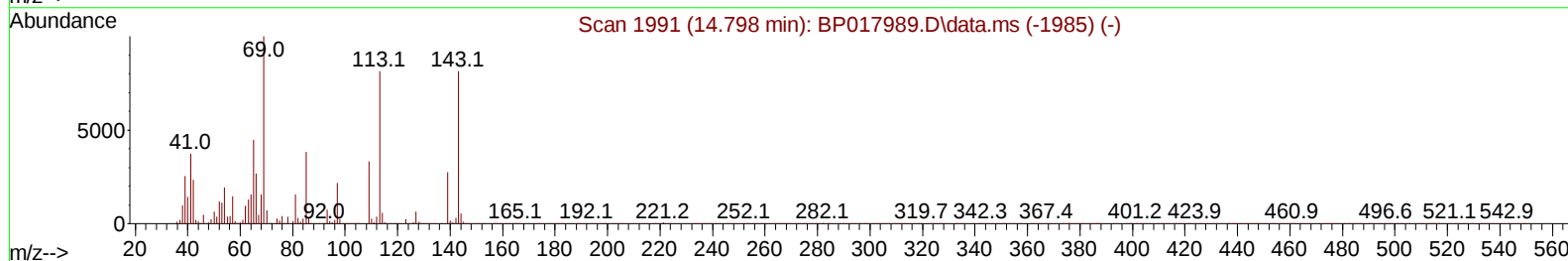
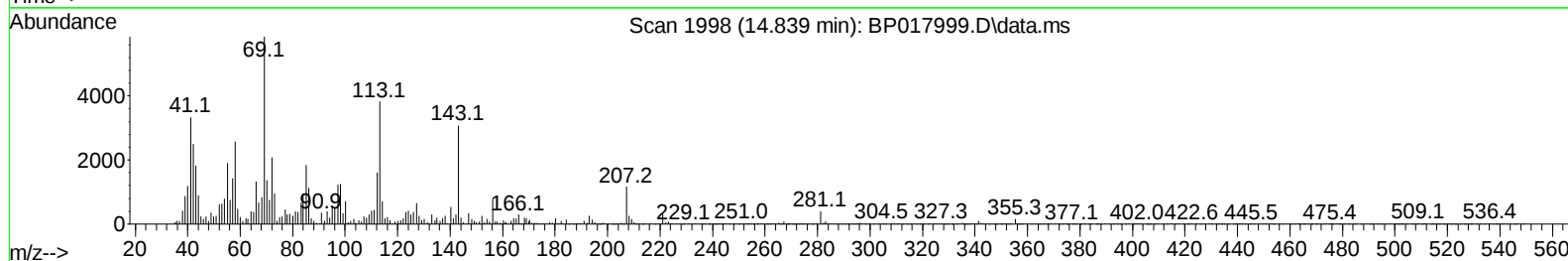
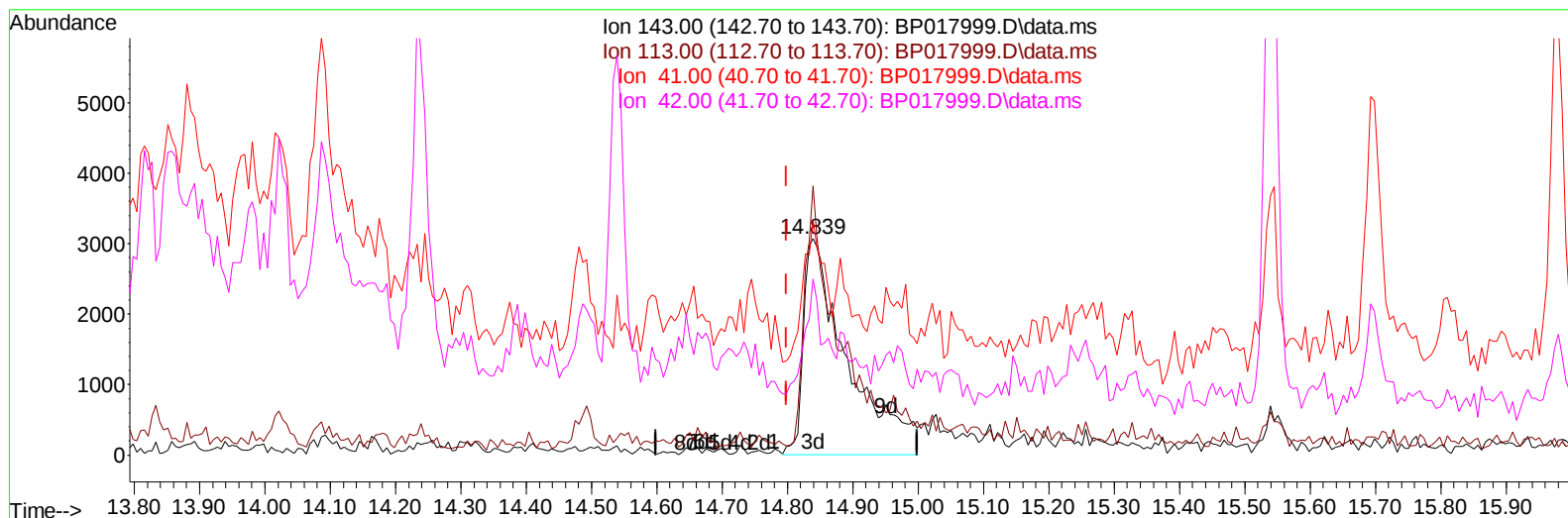
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TIC: BP017999.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.839min (+ 0.041) 6.06 ng/ul m

response 14287

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	100.40	124.13#
41.00	46.90	108.33#
42.00	29.70	81.24#

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Reviewed By :Yogesh Patel 11/11/2023
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Quant Time: Nov 09 23:28:57 2023
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 Quant Title : SVOA CALIBRATION
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.869	152	61764	20.000	ng/ul	0.00
20) Naphthalene-d8	10.686	136	257405	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.539	164	175118	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.321	188	403490	20.000	ng/ul	0.00
79) Chrysene-d12	21.445	240	411399	20.000	ng/ul	0.00
88) Perylene-d12	23.956	264	453150	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.246	96	6941	3.847	ng/uL	0.00
4) Pyridine-d5	3.687	84	62758	13.383	ng/ul	0.00
7) Phenol-d5	7.051	99	49754	8.572	ng/ul	0.02
9) Bis-(2-Chloroethyl)eth...	7.216	67	83676	23.799	ng/ul	0.00
11) 2-Chlorophenol-d4	7.398	132	102203	22.007	ng/ul	0.01
15) 4-Methylphenol-d8	8.604	113	79746	17.130	ng/ul	0.01
21) Nitrobenzene-d5	9.075	128	55851	24.229	ng/ul	0.00
24) 2-Nitrophenol-d4	9.786	143	64758	25.243	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.334	165	105751	23.010	ng/ul	0.02
31) 4-Chloroaniline-d4	10.863	131	169959	26.986	ng/ul	0.00
46) Dimethylphthalate-d6	13.963	166	399027	24.787	ng/ul	0.00
49) Acenaphthylene-d8	14.239	160	413648	23.968	ng/ul	0.00
54) 4-Nitrophenol-d4	14.839	143	14287m	6.060	ng/ul	0.04
60) Fluorene-d10	15.539	176	330940	24.803	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.698	200	64539	22.907	ng/ul	0.00
73) Anthracene-d10	17.421	188	545797	25.245	ng/ul	0.00
81) Pyrene-d10	19.674	212	698895	26.396	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.792	264	680531	25.837	ng/ul	0.00

Target Compounds Qvalue

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

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