

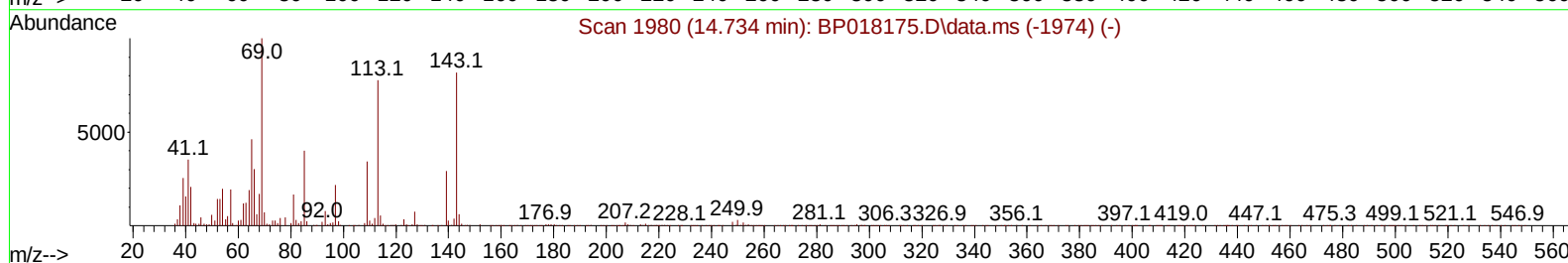
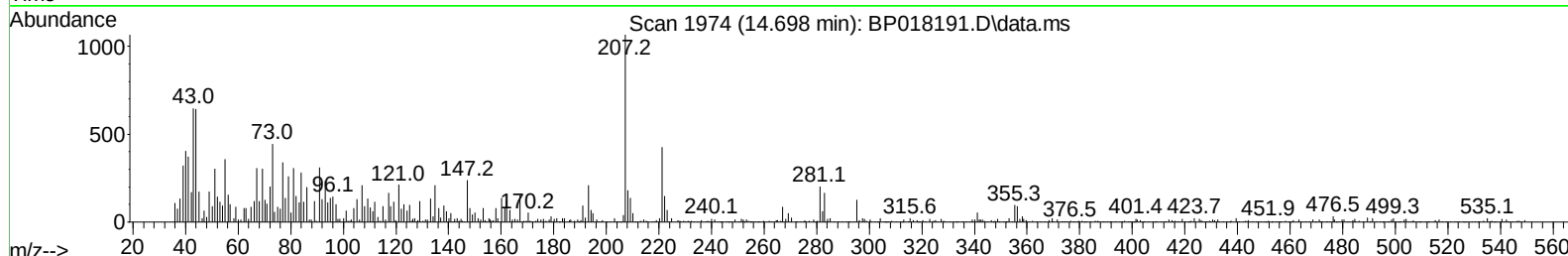
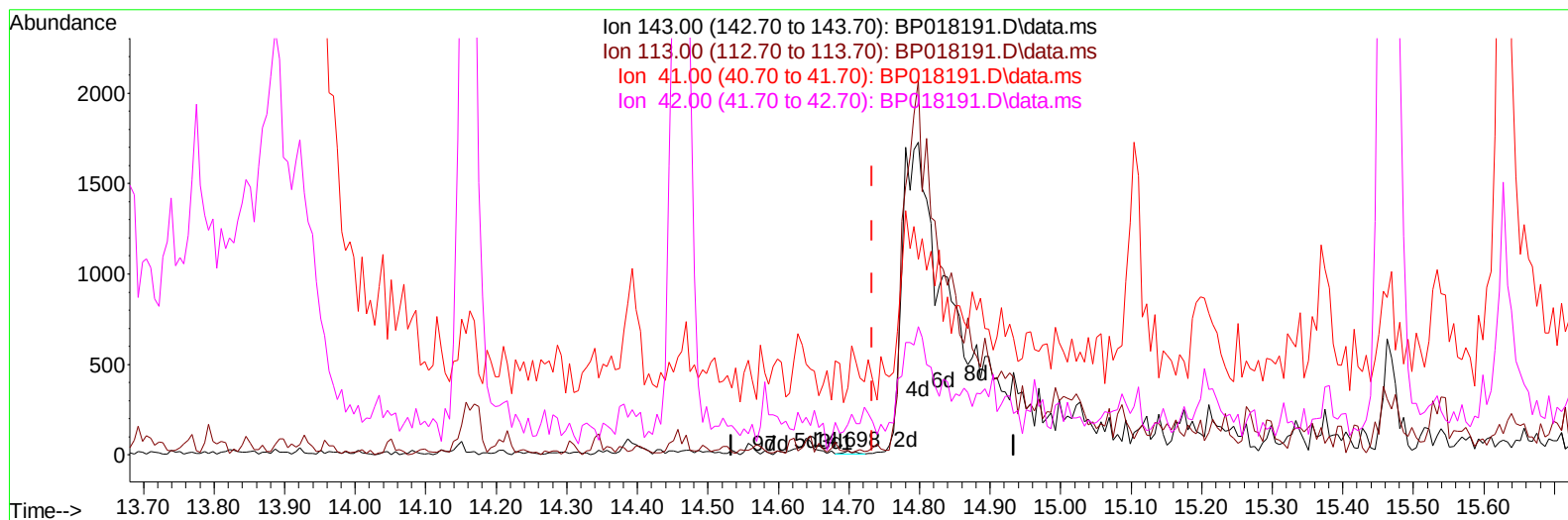
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111523\
 Data File : BP018191.D
 Acq On : 16 Nov 2023 01:35
 Operator : MA/JU
 Sample : 05338-11
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 GCDL6

Manual IntegrationsAPPROVED

Quant Time: Nov 16 02:08:24 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP111023.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Nov 15 18:13:37 2023
 Response via : Initial Calibration

Reviewed By :Yogesh Patel 11/16/2023
 Supervised By :mohammad ahmed 11/17/2023



TIC: BP018191.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.698min (-0.035) 0.01 ng/ul

response 18

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	103.00	133.33#
41.00	47.70	1780.95#
42.00	30.10	804.76#

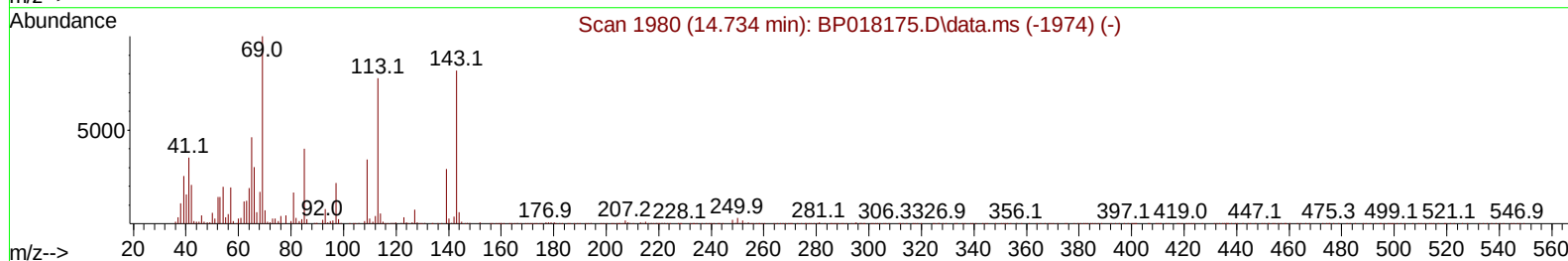
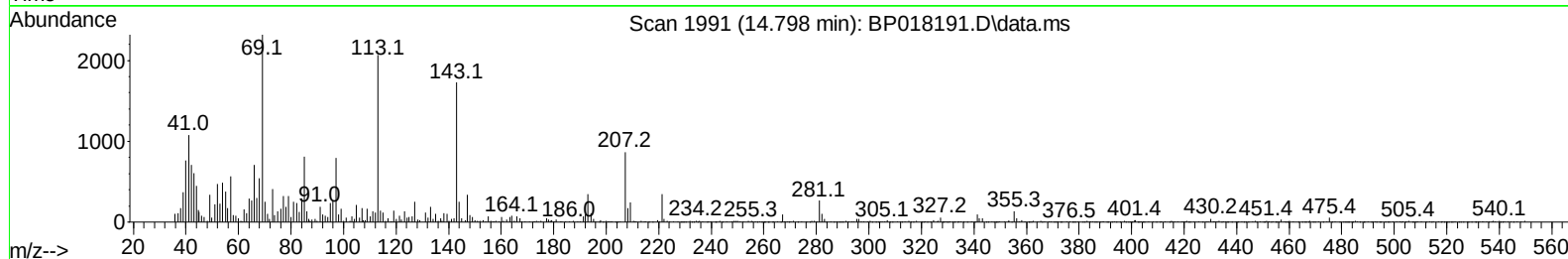
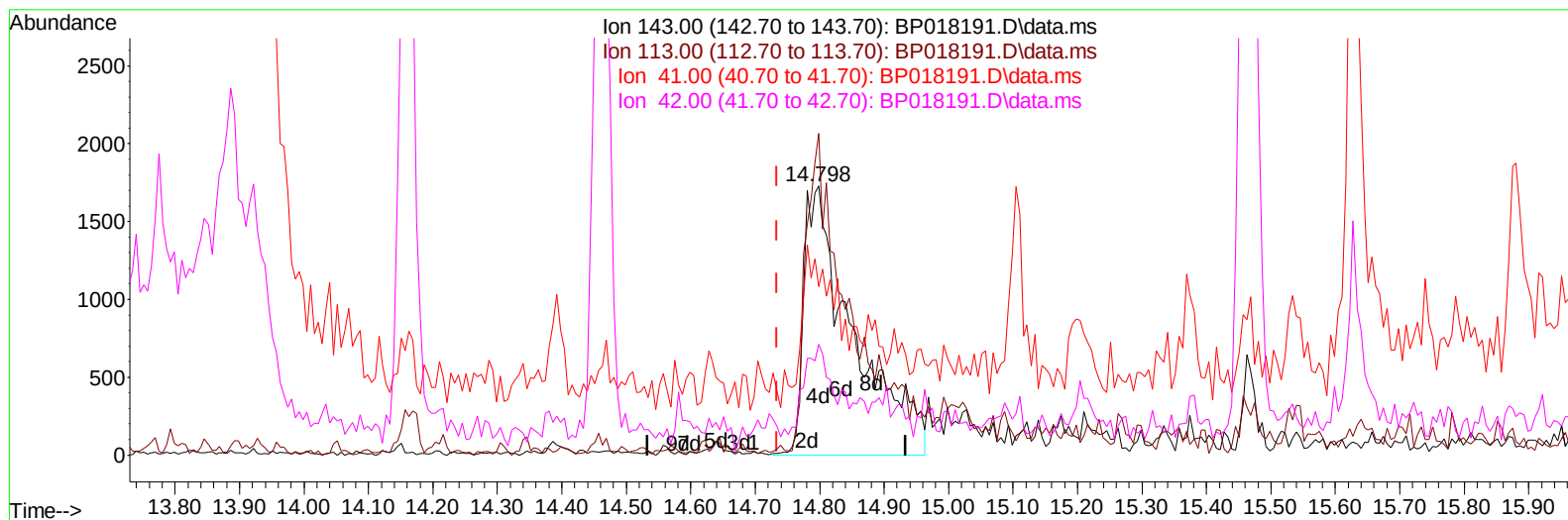
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(54) 4-Nitrophenol-d4 (S)

14.798min (+ 0.065) 3.98 ng/ul m

response 9179

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	103.00	119.62
41.00	47.70	62.50#
42.00	30.10	41.03#

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.787	152	71387	20.000	ng/ul	0.00
20) Naphthalene-d8	10.598	136	274973	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.463	164	167227	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.245	188	391561	20.000	ng/ul	0.00
79) Chrysene-d12	21.374	240	397066	20.000	ng/ul	0.00
88) Perylene-d12	23.845	264	432313	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.164	96	7197	3.556	ng/uL	0.00
4) Pyridine-d5	3.605	84	37811	7.166	ng/ul	0.01
7) Phenol-d5	6.969	99	34461	5.000	ng/ul	0.02
9) Bis-(2-Chloroethyl)eth...	7.134	67	100541	24.824	ng/ul	0.00
11) 2-Chlorophenol-d4	7.305	132	101142	18.532	ng/ul	0.00
15) 4-Methylphenol-d8	8.516	113	67224	11.903	ng/ul	0.00
21) Nitrobenzene-d5	8.993	128	57071	23.037	ng/ul	0.00
24) 2-Nitrophenol-d4	9.704	143	62174	22.153	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.234	165	101177	20.226	ng/ul	0.00
31) 4-Chloroaniline-d4	10.793	131	88228	12.861	ng/ul	0.00
46) Dimethylphthalate-d6	13.887	166	399360	25.883	ng/ul	0.00
49) Acenaphthylene-d8	14.163	160	400662	23.841	ng/ul	0.00
54) 4-Nitrophenol-d4	14.798	143	9179m	3.981	ng/ul	0.06
60) Fluorene-d10	15.463	176	323495	25.394	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.628	200	52109	18.409	ng/ul	0.00
73) Anthracene-d10	17.345	188	549657	25.955	ng/ul	0.00
81) Pyrene-d10	19.598	212	696646	26.266	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.680	264	682019	27.131	ng/ul	0.00

Target Compounds Qvalue

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

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