

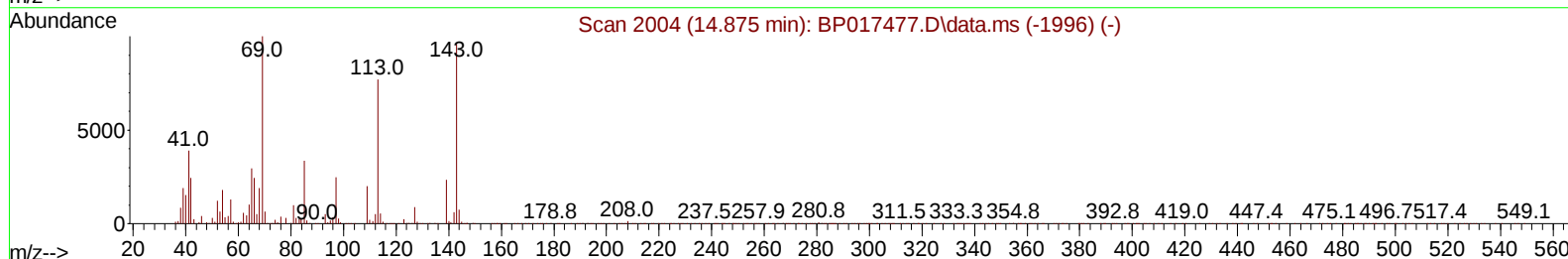
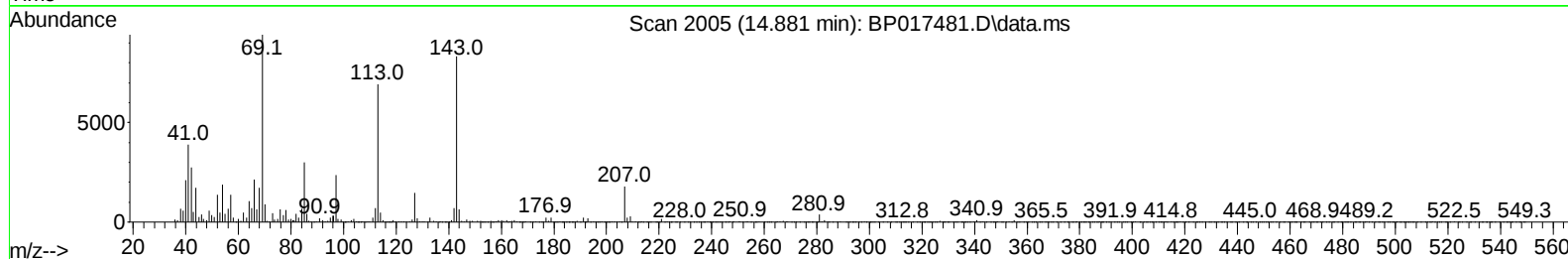
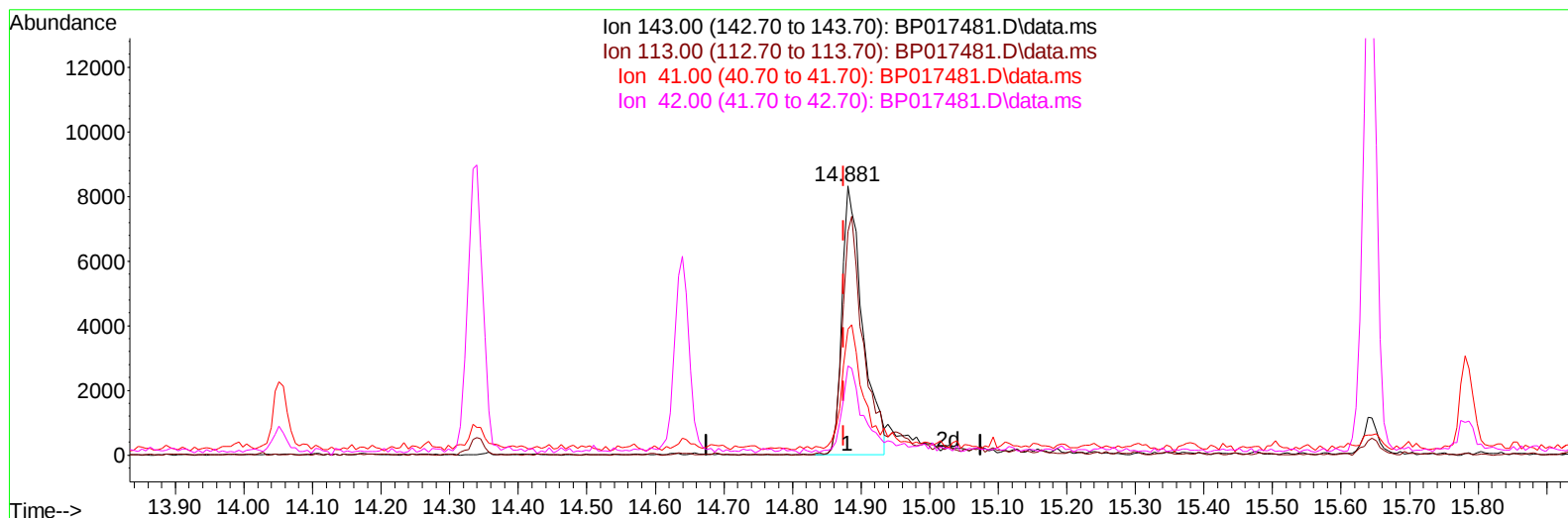
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100223\
 Data File : BP017481.D
 Acq On : 02 Oct 2023 17:24
 Operator : MA/JU
 Sample : 04571-09
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BBJF0

Manual IntegrationsAPPROVED

Quant Time: Oct 02 23:14:56 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092023.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Sat Sep 30 04:02:20 2023
 Response via : Initial Calibration

Reviewed By :Yogesh Patel 10/03/2023
 Supervised By :mohammad ahmed 10/05/2023



TIC: BP017481.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.881min (+ 0.006) 5.41 ng/ul

response 17316

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	74.90	83.06
41.00	36.70	46.72#
42.00	26.40	33.14#

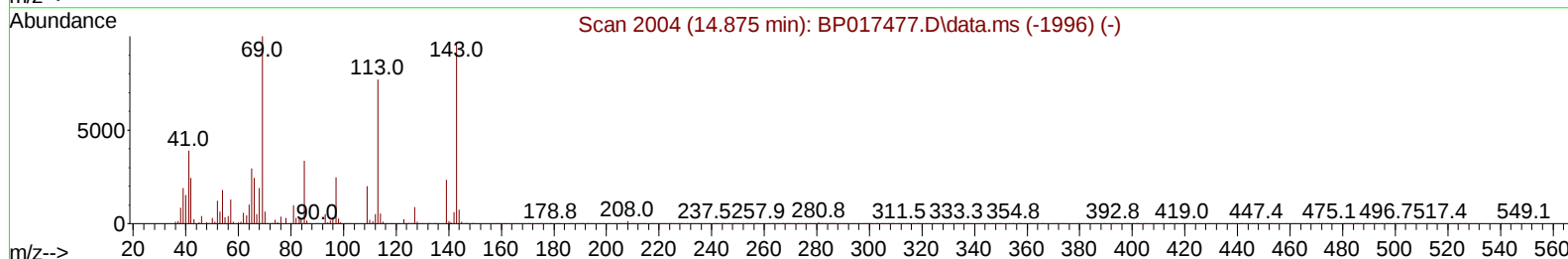
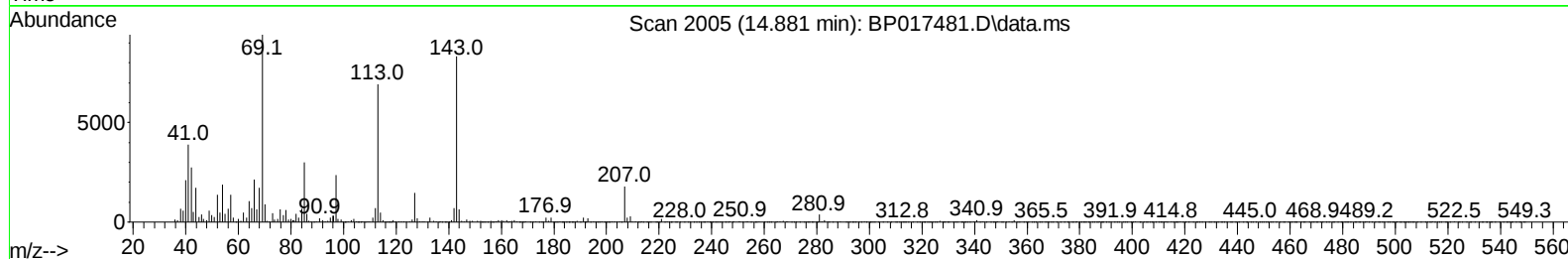
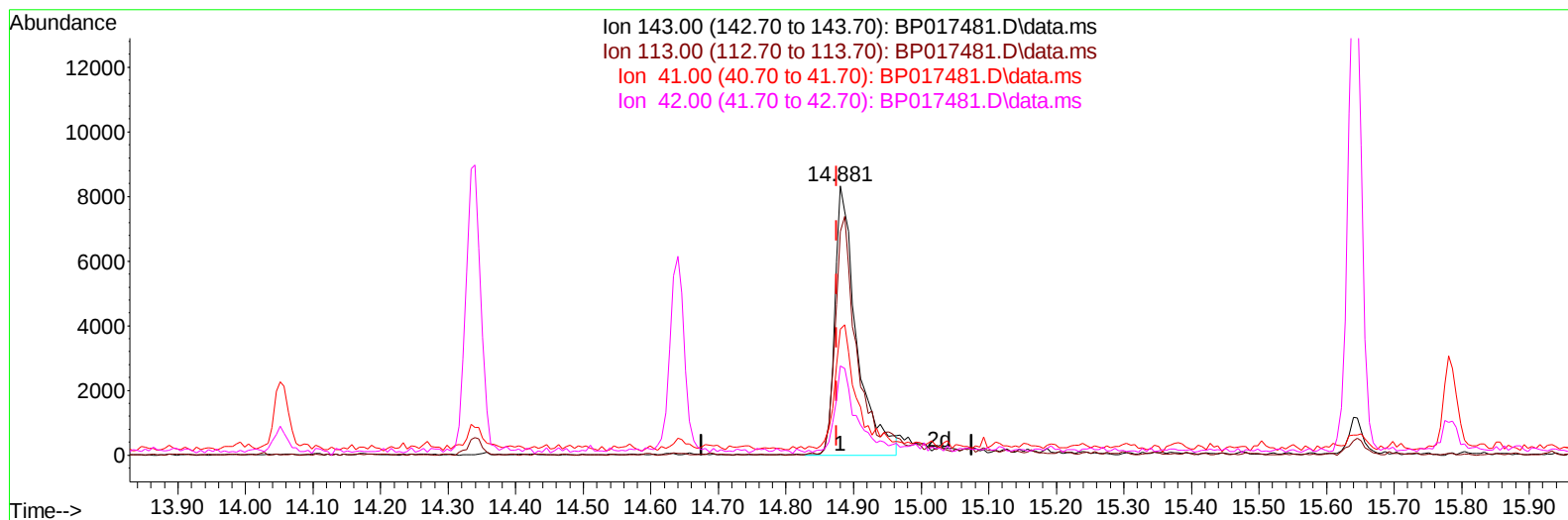
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TIC: BP017481.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.881min (+ 0.006) 5.80 ng/ul m

response 18582

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	74.90	83.06
41.00	36.70	46.72#
42.00	26.40	33.14#

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Instrument :
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Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 10/03/2023
 Supervised By :mohammad ahmed 10/05/2023

Quant Time: Oct 02 23:25:31 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092023.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Sat Sep 30 04:02:20 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.946	152	114478	20.000	ng/ul	0.00
20) Naphthalene-d8	10.787	136	442765	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.639	164	256695	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.428	188	550794	20.000	ng/ul	0.00
79) Chrysene-d12	21.557	240	533540	20.000	ng/ul	0.00
88) Perylene-d12	24.121	264	599851	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.293	96	13811	4.955	ng/uL	0.00
4) Pyridine-d5	3.734	84	73565	9.442	ng/ul	0.00
7) Phenol-d5	7.105	99	115680	12.304	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.287	67	230800	40.676	ng/ul	0.00
11) 2-Chlorophenol-d4	7.463	132	157002	21.054	ng/ul	0.00
15) 4-Methylphenol-d8	8.669	113	136022	18.609	ng/ul	0.00
21) Nitrobenzene-d5	9.151	128	146430	45.746	ng/ul	0.00
24) 2-Nitrophenol-d4	9.869	143	85488	24.597	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.398	165	156812	23.881	ng/ul	0.00
31) 4-Chloroaniline-d4	10.957	131	102913	10.819	ng/ul	0.00
46) Dimethylphthalate-d6	14.057	166	907719	49.362	ng/ul	0.00
49) Acenaphthylene-d8	14.339	160	975491	46.124	ng/ul	0.00
54) 4-Nitrophenol-d4	14.881	143	18582m	5.800	ng/ul	0.00
60) Fluorene-d10	15.645	176	786942	49.709	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.786	200	45125	14.536	ng/ul	0.00
73) Anthracene-d10	17.528	188	1308647	53.571	ng/ul	0.00
81) Pyrene-d10	19.780	212	1588752	55.065	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.957	264	1503304	52.052	ng/ul	0.00

Target Compounds Qvalue

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

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