

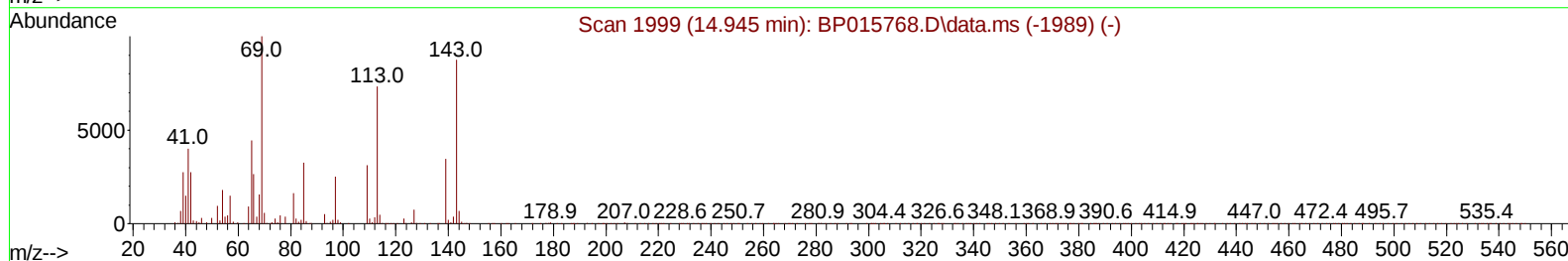
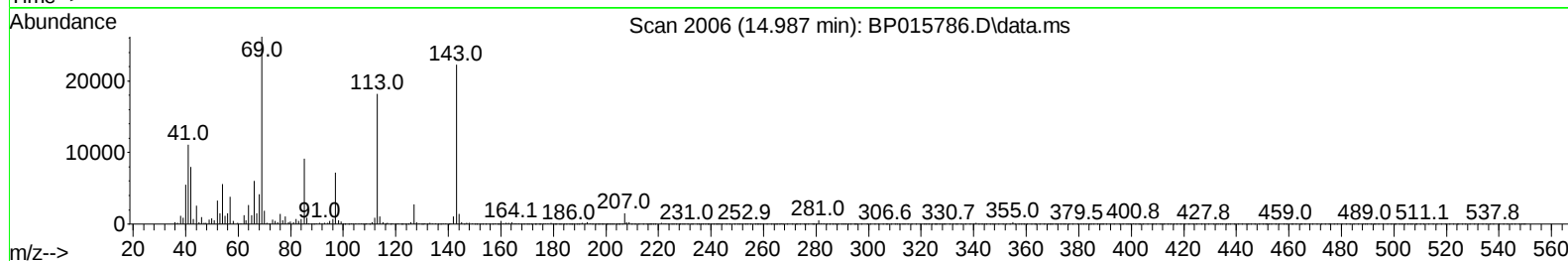
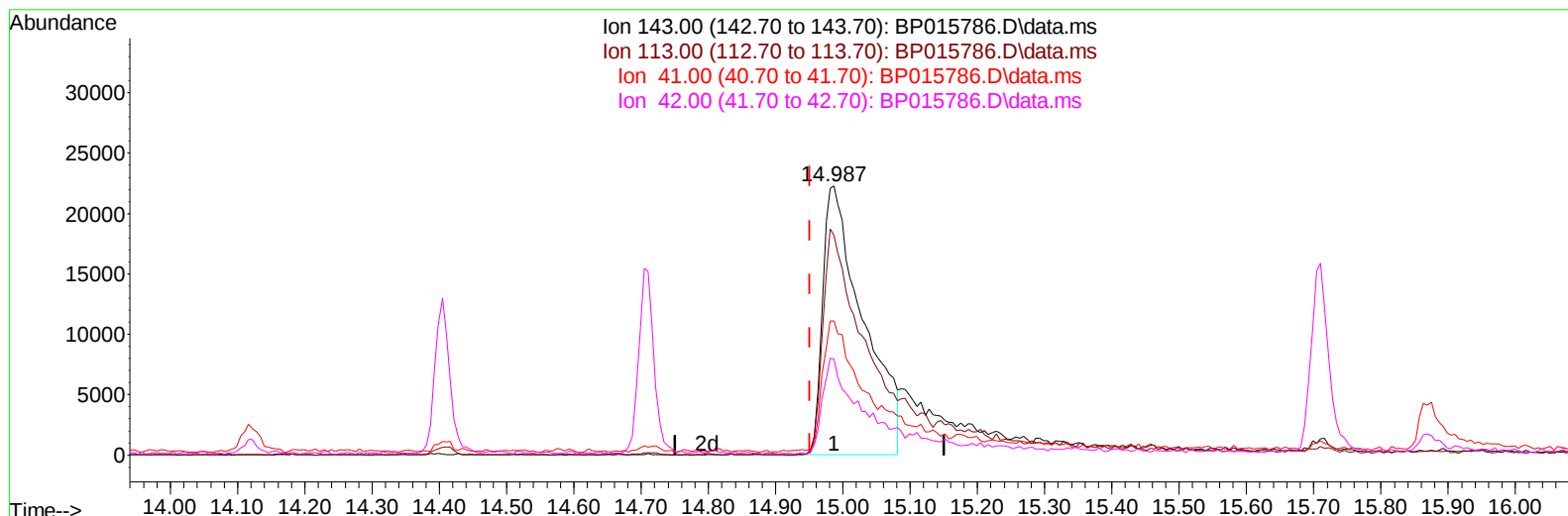
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062823\
 Data File : BP015786.D
 Acq On : 28 Jun 2023 19:30
 Operator : MA/JU
 Sample : PB153625BL
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK625

Manual IntegrationsAPPROVED

Quant Time: Jun 29 02:31:32 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062623.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Jun 29 02:30:23 2023
 Response via : Initial Calibration

Reviewed By :Yogesh Patel 06/29/2023
 Supervised By :mohammad ahmed 06/29/2023



TIC: BP015786.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.987min (+ 0.035) 14.32 ng/u1

response 92509

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	94.00	81.75
41.00	45.70	49.92
42.00	31.20	35.75

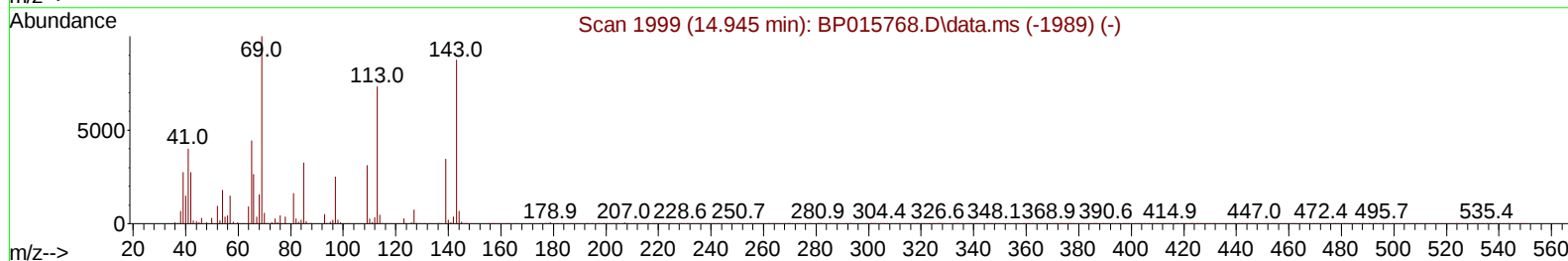
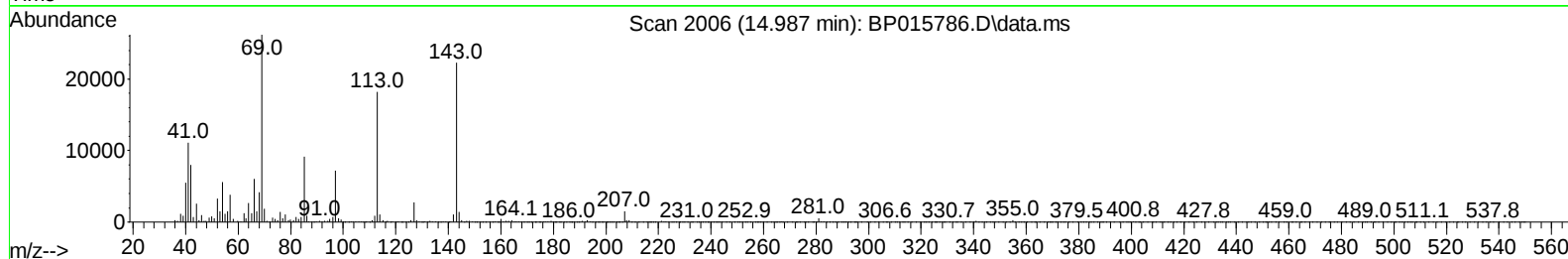
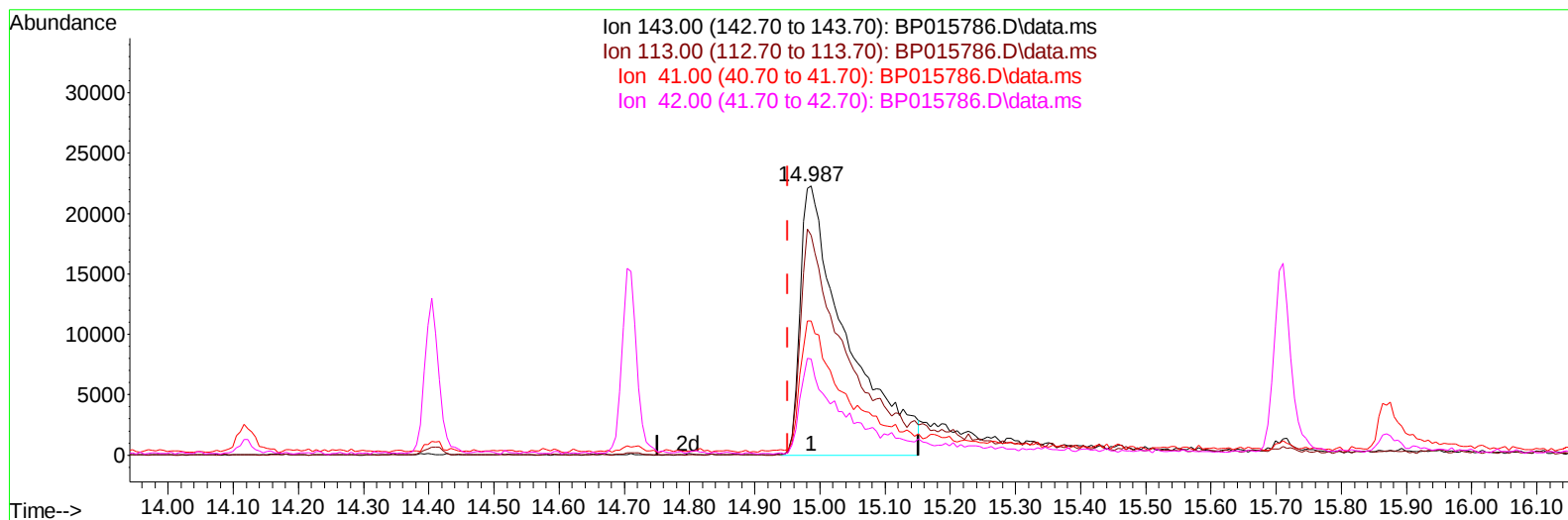
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TIC: BP015786.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.987min (+ 0.035) 16.99 ng/ul m

response 109744

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	94.00	81.75
41.00	45.70	49.92
42.00	31.20	35.75

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.058	152	269344	20.000	ng/ul	0.00
20) Naphthalene-d8	10.887	136	1118033	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.710	164	633464	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.469	188	1156633	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.563	240	918614	20.000	ng/ul	0.00
88) Perylene-d12	24.104	264	1101865	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.381	96	31187	4.166	ng/uL	0.00
4) Pyridine-d5	3.823	84	395445	20.955	ng/ul	0.00
7) Phenol-d5	7.228	99	513194	22.269	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.381	67	350817	24.605	ng/ul	0.00
11) 2-Chlorophenol-d4	7.587	132	415253	23.870	ng/ul	0.00
15) 4-Methylphenol-d8	8.787	113	425799	23.388	ng/ul	0.00
21) Nitrobenzene-d5	9.246	128	204770	23.965	ng/ul	0.00
24) 2-Nitrophenol-d4	9.975	143	231561	23.220	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.528	165	368004	20.708	ng/ul	0.01
31) 4-Chloroaniline-d4	11.046	131	517230	22.657	ng/ul	0.01
46) Dimethylphthalate-d6	14.116	166	1201837	24.546	ng/ul	0.00
49) Acenaphthylene-d8	14.404	160	1341847	24.380	ng/ul	0.00
54) 4-Nitrophenol-d4	14.987	143	109744m	16.985	ng/ul	0.04
60) Fluorene-d10	15.710	176	912542	22.635	ng/ul	0.01
65) 4,6-Dinitro-2-methylph...	15.869	200	112254	15.781	ng/ul	0.01
73) Anthracene-d10	17.575	188	1299645	24.599	ng/ul	0.00
81) Pyrene-d10	19.798	212	1352766	22.552	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.933	264	1396895	25.132	ng/ul	0.00

Target Compounds Qvalue

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

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