

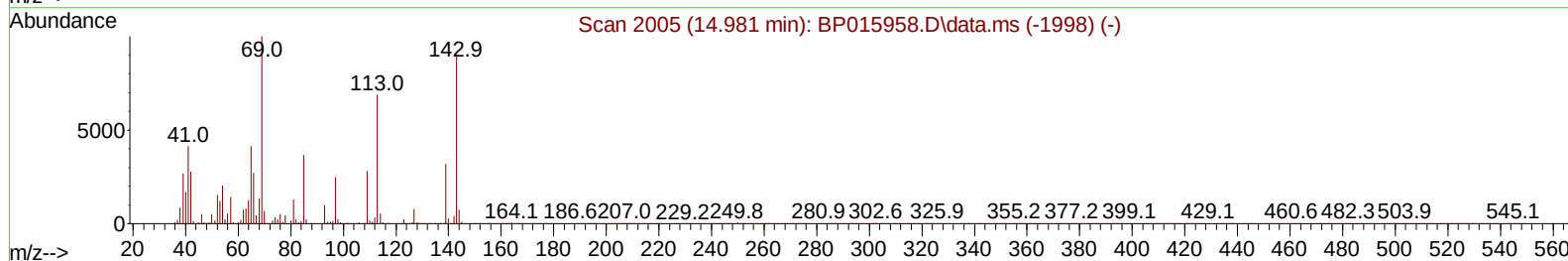
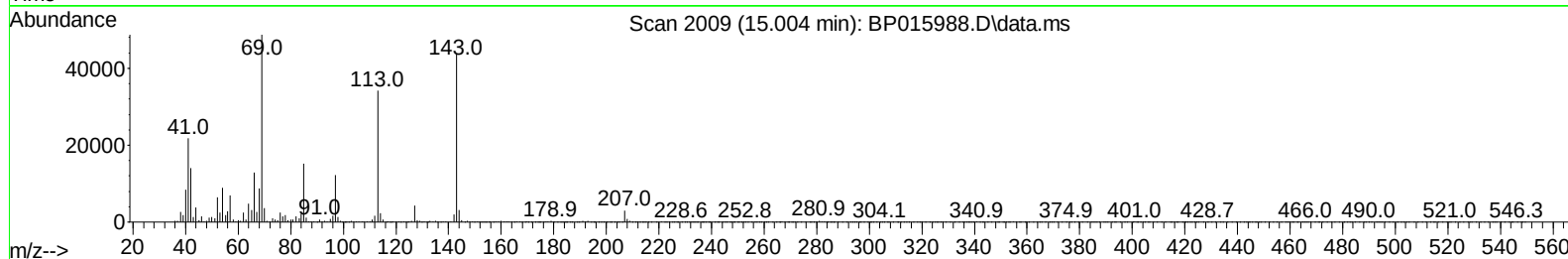
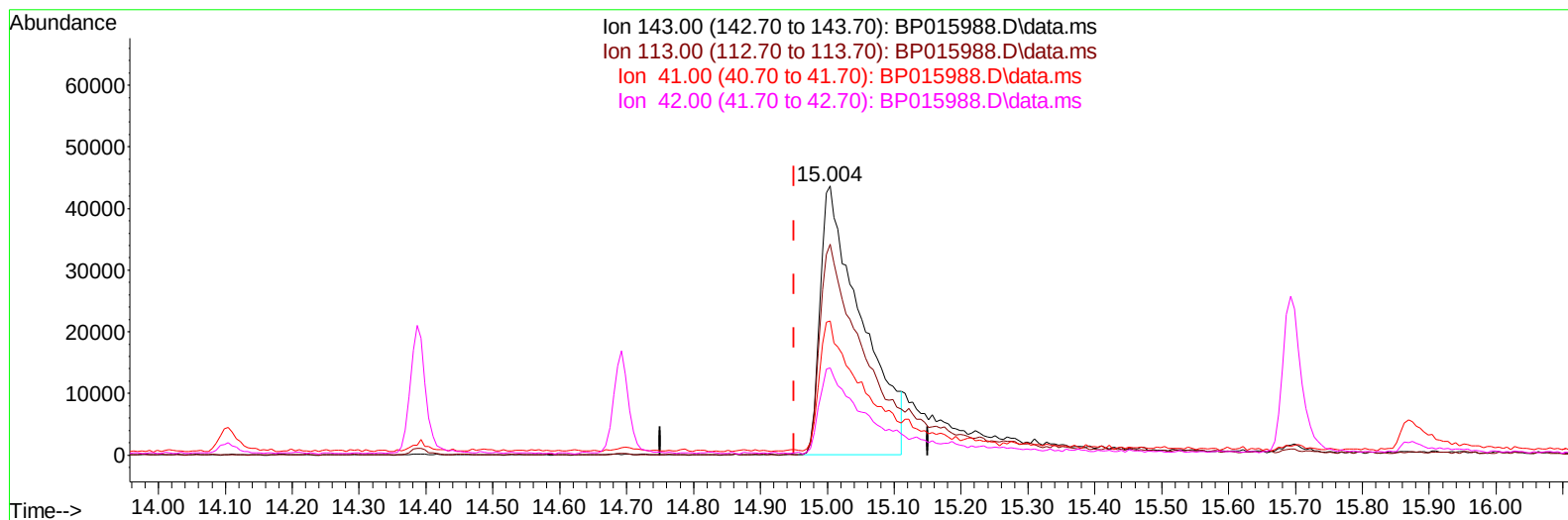
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP070423\
Data File : BP015988.D
Acq On : 04 Jul 2023 22:29
Operator : MA/JU
Sample : PB153766BL
Misc :
ALS Vial : 35 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
SBLK766

Manual IntegrationsAPPROVED

Quant Time: Jul 04 23:45:03 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062623.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Sun Jul 02 06:12:16 2023
Response via : Initial Calibration

Reviewed By :Yogesh Patel 07/05/2023
Supervised By :mohammad ahmed 07/05/2023



TIC: BP015988.D\data.ms

(54) 4-Nitrophenol-d4 (S)

15.004min (+ 0.053) 26.15 ng/ul

response 185951

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	94.00	78.31
41.00	45.70	49.94
42.00	31.20	32.36

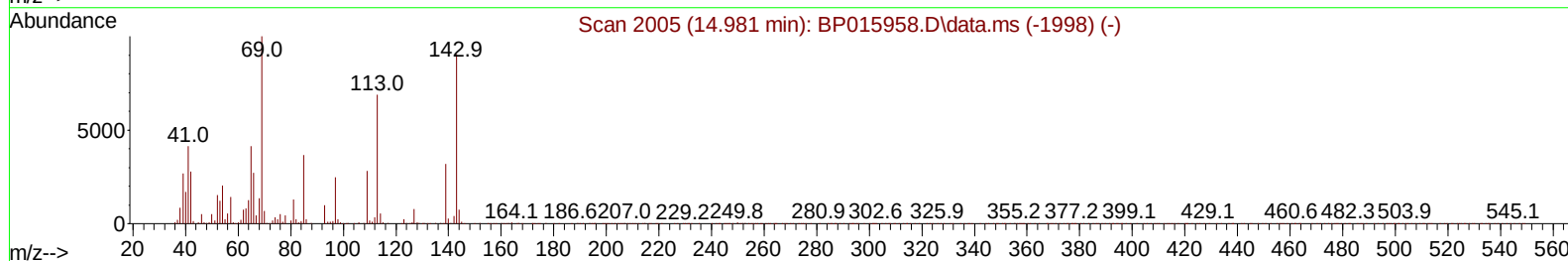
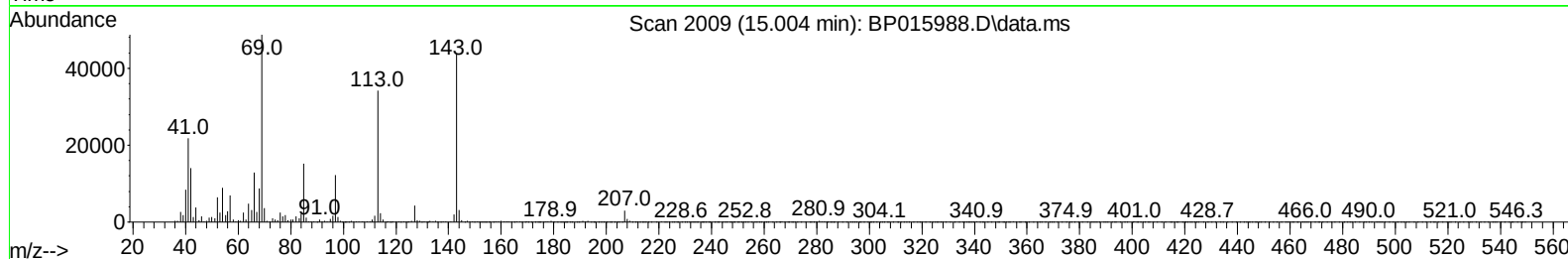
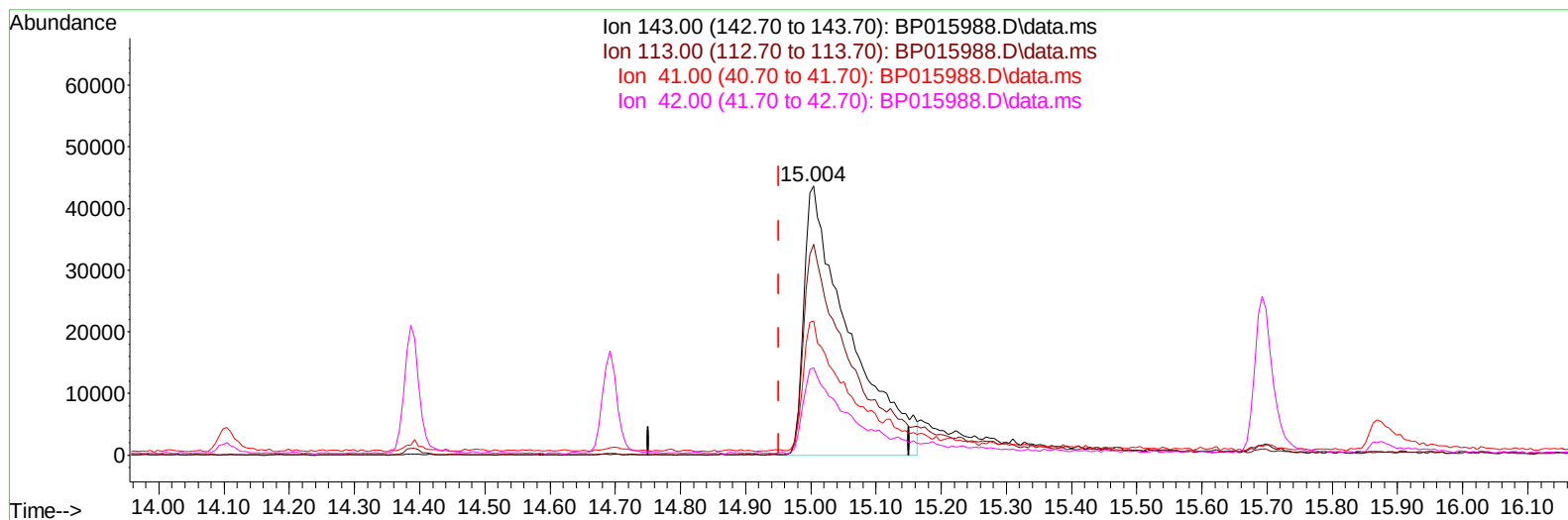
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TIC: BP015988.D\data.ms

(54) 4-Nitrophenol-d4 (S)

15.004min (+ 0.053) 29.46 ng/ul m

response 209493

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	94.00	78.31
41.00	45.70	49.94
42.00	31.20	32.36

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Instrument :
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Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 07/05/2023
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Quant Time: Jul 04 23:45:47 2023
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.040	152	254445	20.000	ng/ul	0.00
20) Naphthalene-d8	10.869	136	1164280	20.000	ng/ul	# 0.00
38) Acenaphthene-d10	14.693	164	697145	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.457	188	1335703	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.557	240	938536	20.000	ng/ul	# 0.00
88) Perylene-d12	24.092	264	1030539	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.370	96	46110	6.520	ng/uL	0.00
4) Pyridine-d5	3.811	84	635618	35.654	ng/ul	0.00
7) Phenol-d5	7.228	99	800735	36.780	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.363	67	556174	41.293	ng/ul	0.00
11) 2-Chlorophenol-d4	7.575	132	652411	39.699	ng/ul	0.00
15) 4-Methylphenol-d8	8.781	113	670292	38.974	ng/ul	0.00
21) Nitrobenzene-d5	9.228	128	326285	36.670	ng/ul	0.00
24) 2-Nitrophenol-d4	9.963	143	358223	34.495	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.534	165	583287	31.519	ng/ul	0.02
31) 4-Chloroaniline-d4	11.040	131	857680	36.079	ng/ul	0.01
46) Dimethylphthalate-d6	14.104	166	1977929	36.707	ng/ul	0.00
49) Acenaphthylene-d8	14.387	160	2203056	36.370	ng/ul	0.00
54) 4-Nitrophenol-d4	15.004	143	209493m	29.461	ng/ul	0.05
60) Fluorene-d10	15.692	176	1605023	36.175	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.875	200	147601	17.968	ng/ul	0.02
73) Anthracene-d10	17.557	188	2355491	38.607	ng/ul	0.00
81) Pyrene-d10	19.786	212	2363656	38.569	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.927	264	1998742	38.448	ng/ul	0.00

Target Compounds Qvalue

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

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