

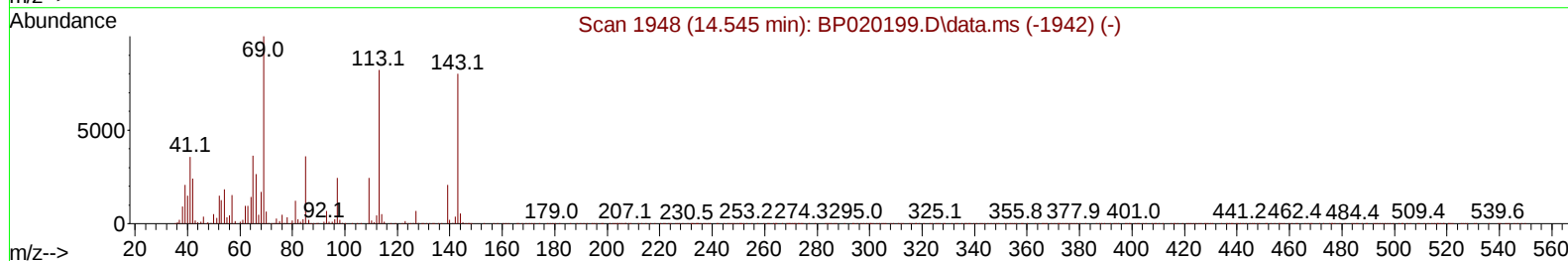
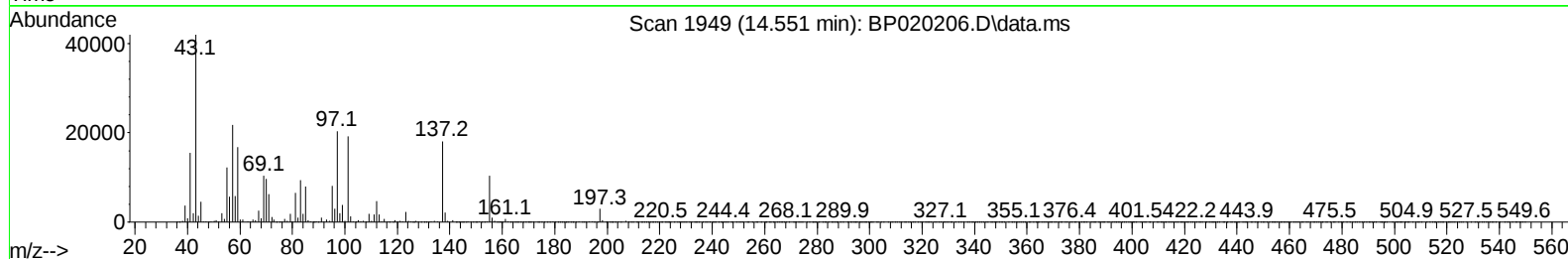
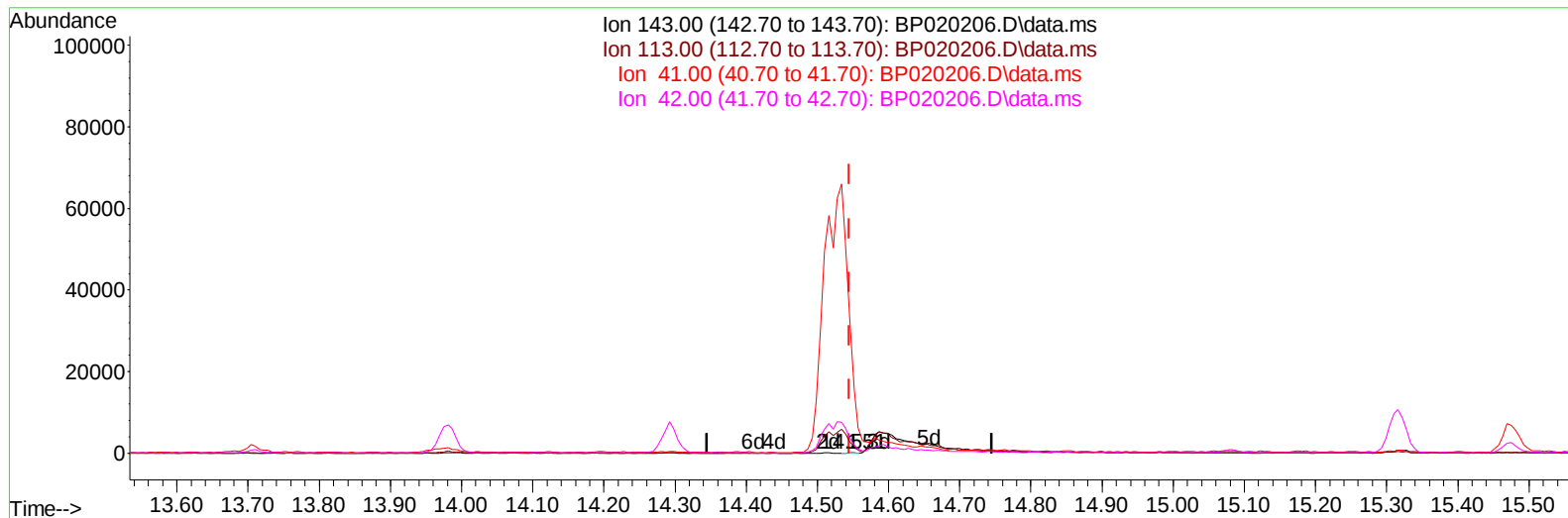
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP050624\
 Data File : BP020206.D
 Acq On : 06 May 2024 14:22
 Operator : MA/JU
 Sample : P2279-06
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0RQ9

Manual IntegrationsAPPROVED

Quant Time: May 06 15:28:05 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP050124.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon May 06 12:26:12 2024
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 05/07/2024
 Supervised By :mohammad ahmed 05/08/2024



TIC: BP020206.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.551min (+ 0.006) 0.01 ng/u1

response 22

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	101.40	5917.86#
41.00	50.30	55510.71#
42.00	32.50	6982.14#

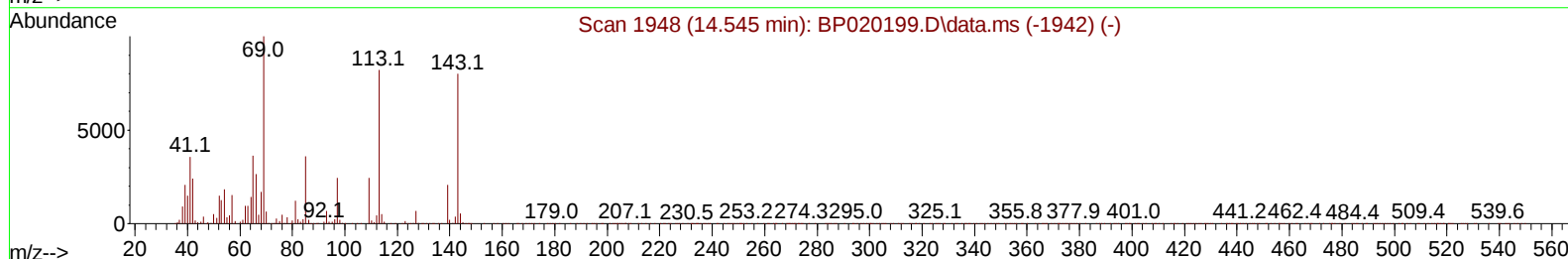
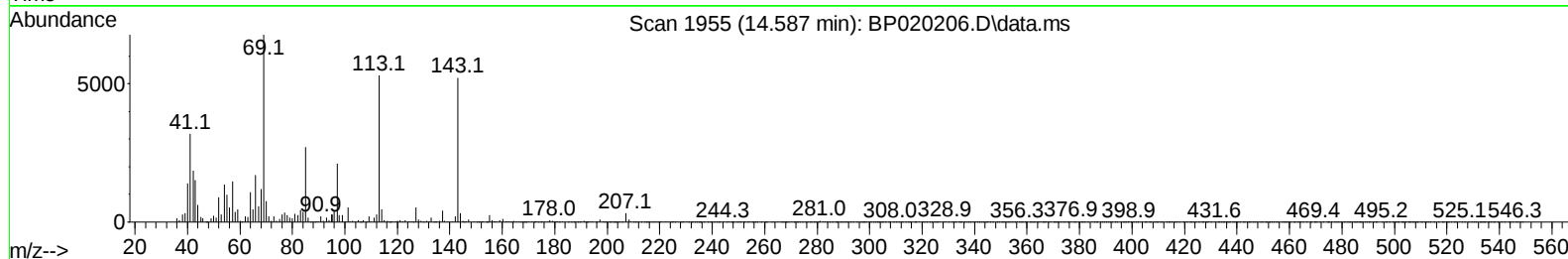
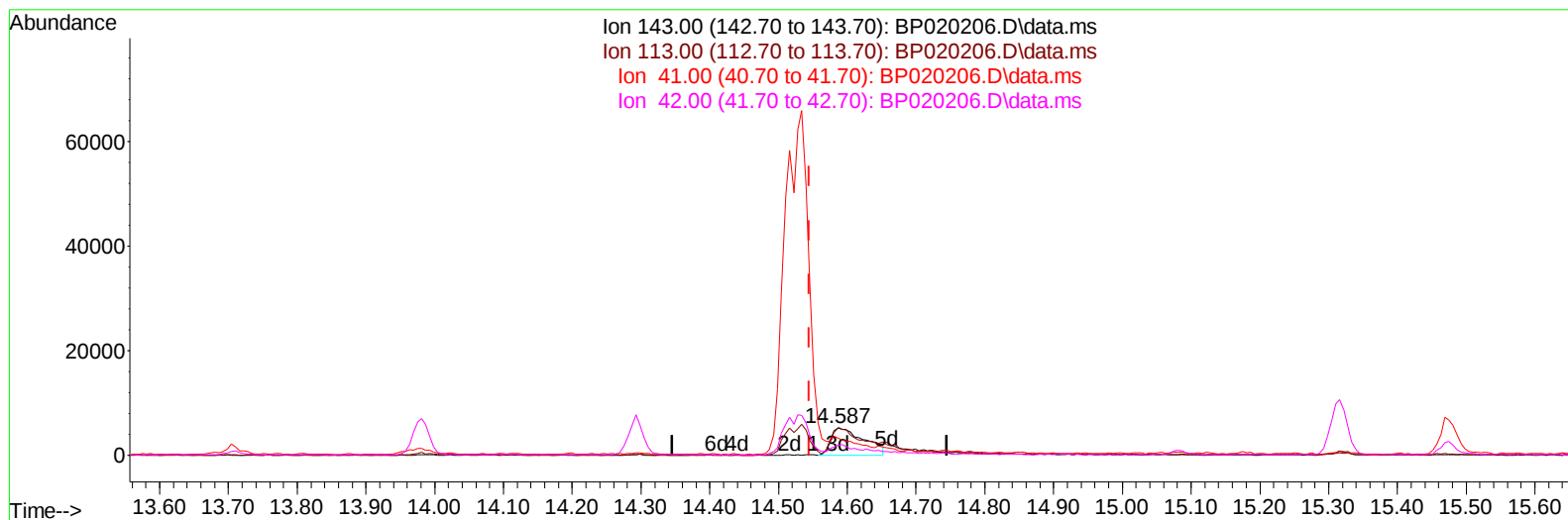
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(54) 4-Nitrophenol-d4 (S)

14.587min (+ 0.041) 5.25 ng/ul m

response 17550

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	101.40	101.59
41.00	50.30	61.11#
42.00	32.50	35.60

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 Misc :
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Instrument :
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Manual IntegrationsAPPROVED

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.622	152	95579	20.000	ng/ul	0.02
20) Naphthalene-d8	10.393	136	405050	20.000	ng/ul	0.02
38) Acenaphthene-d10	14.293	164	297847	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.140	188	670267	20.000	ng/ul	0.00
79) Chrysene-d12	21.622	240	658044	20.000	ng/ul	-0.01
88) Perylene-d12	24.969	264	627103	20.000	ng/ul	-0.03
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.217	96	9306	4.112	ng/uL	0.00
4) Pyridine-d5	3.634	84	38300	5.824	ng/ul	0.03
7) Phenol-d5	6.822	99	48855	5.952	ng/ul	0.04
9) Bis-(2-Chloroethyl)eth...	6.975	67	142337	28.546	ng/ul	0.02
11) 2-Chlorophenol-d4	7.164	132	137064	23.358	ng/ul	0.02
15) 4-Methylphenol-d8	8.334	113	98447	14.835	ng/ul	0.02
21) Nitrobenzene-d5	8.781	128	90965	29.990	ng/ul	0.02
24) 2-Nitrophenol-d4	9.493	143	101267	29.155	ng/ul	0.02
28) 2,4-Dichlorophenol-d3	10.034	165	173786	27.418	ng/ul	0.02
31) 4-Chloroaniline-d4	10.558	131	182955	20.624	ng/ul	0.02
46) Dimethylphthalate-d6	13.710	166	722373	33.212	ng/ul	0.00
49) Acenaphthylene-d8	13.981	160	736869	30.668	ng/ul	0.00
54) 4-Nitrophenol-d4	14.587	143	17550m	5.250	ng/ul	0.04
60) Fluorene-d10	15.316	176	614434	33.868	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.475	200	135851	31.940	ng/ul	0.00
73) Anthracene-d10	17.240	188	1070860	36.931	ng/ul	0.00
81) Pyrene-d10	19.639	212	1354718	35.219	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.757	264	1130819	37.295	ng/ul	0.00

Target Compounds Qvalue

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

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