

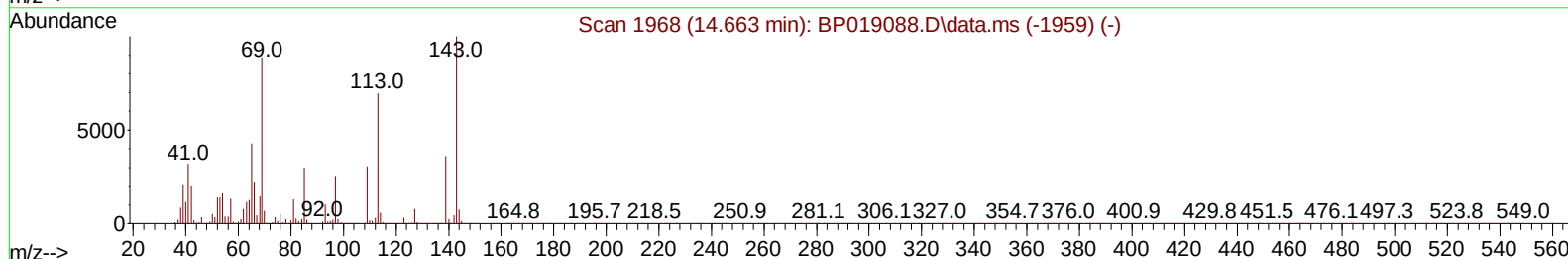
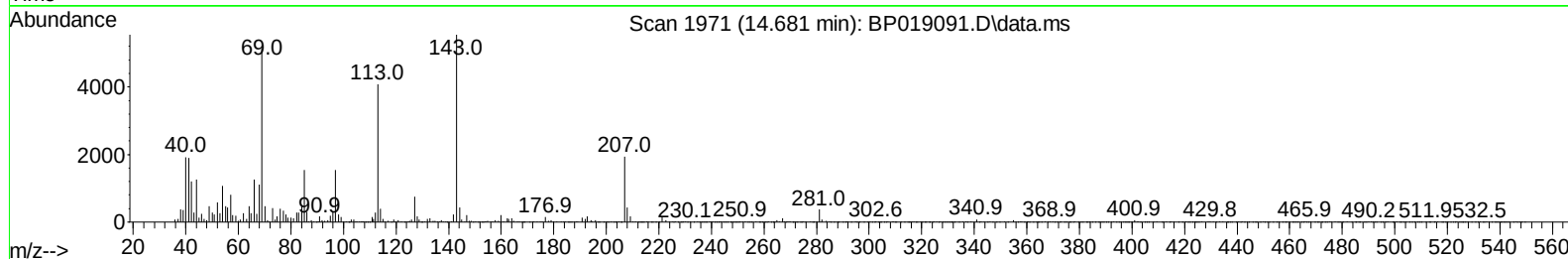
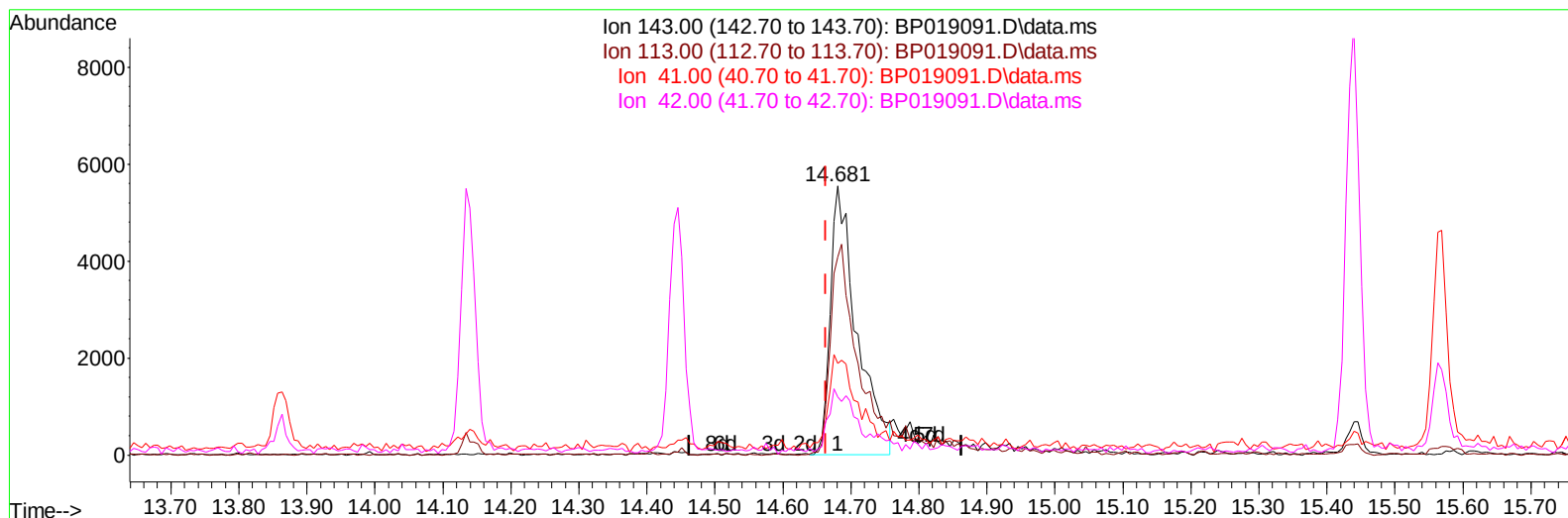
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122623\
 Data File : BP019091.D
 Acq On : 26 Dec 2023 20:02
 Operator : MA/JU
 Sample : 05976-03
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BH7W0

Manual IntegrationsAPPROVED

Quant Time: Dec 26 22:20:57 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120123.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Dec 26 22:20:53 2023
 Response via : Initial Calibration

Reviewed By :Yogesh Patel 12/28/2023
 Supervised By :mohammad ahmed 12/28/2023



TIC: BP019091.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.681min (+ 0.018) 3.41 ng/ul

response 15054

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	70.80	73.66
41.00	34.60	34.13
42.00	23.00	21.71

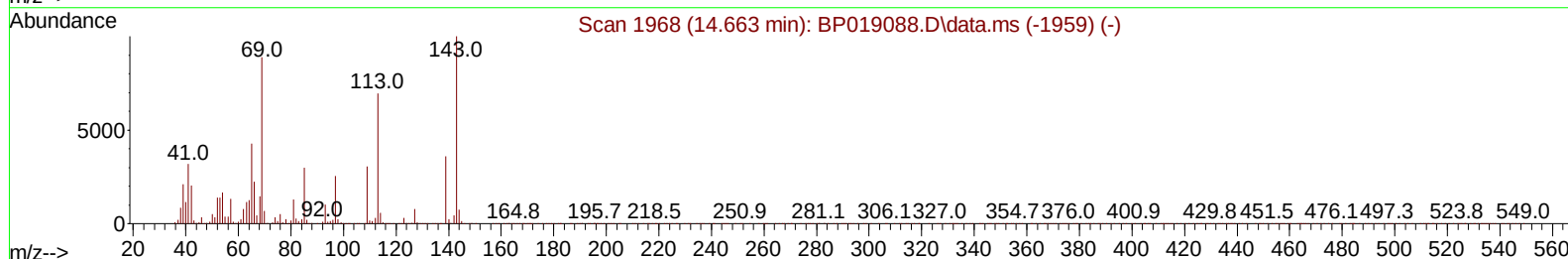
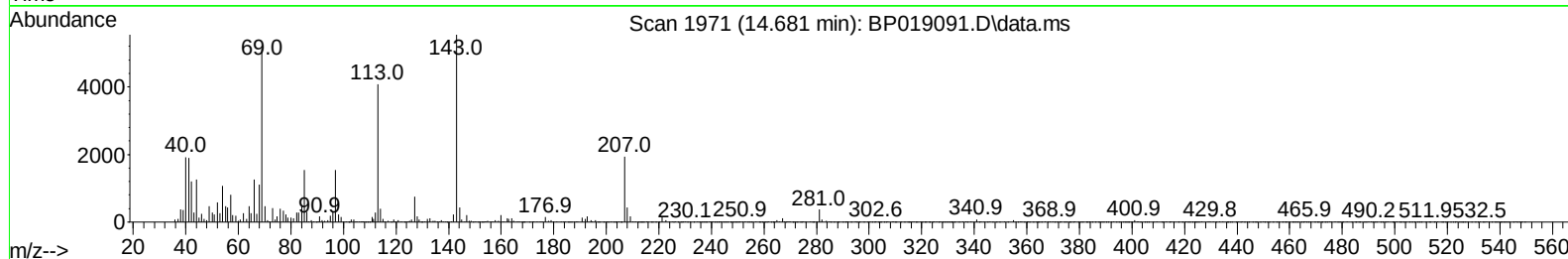
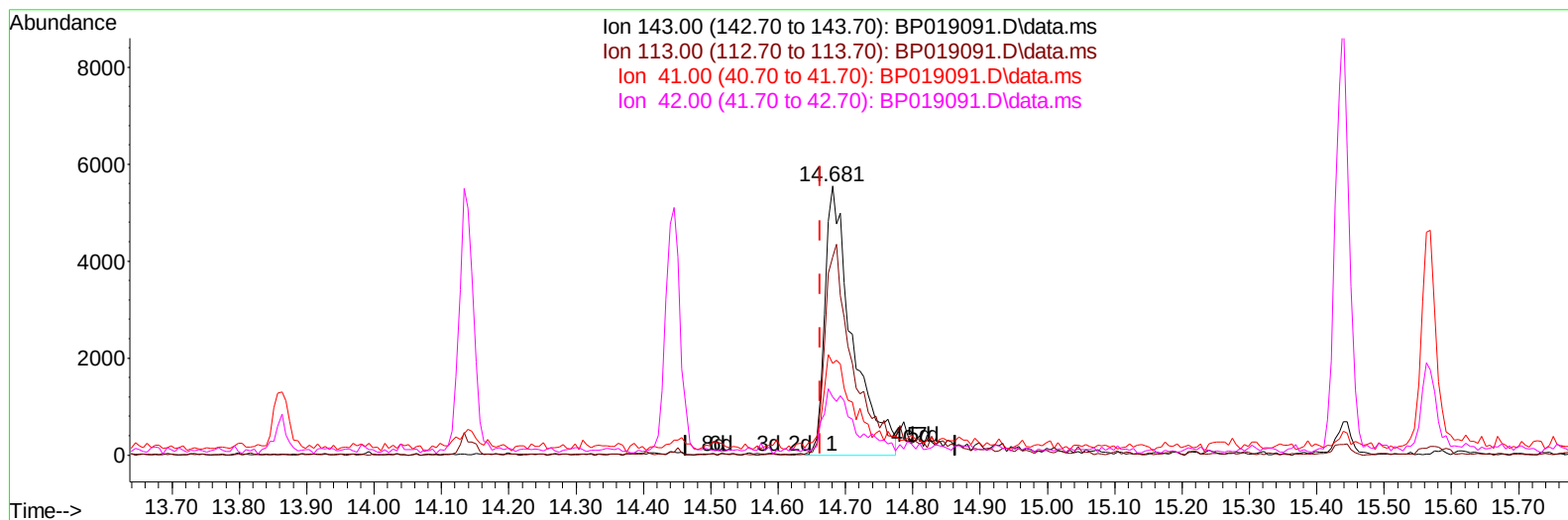
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(54) 4-Nitrophenol-d4 (S)

14.681min (+ 0.018) 3.55 ng/ul m

response 15660

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	70.80	73.66
41.00	34.60	34.13
42.00	23.00	21.71

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.810	152	134140	20.000	ng/ul	0.00
20) Naphthalene-d8	10.604	136	495154	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.445	164	301295	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.198	188	674563	20.000	ng/ul	0.00
79) Chrysene-d12	21.286	240	732332	20.000	ng/ul	0.00
88) Perylene-d12	23.645	264	924526	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.246	96	14576	3.703	ng/uL	0.00
4) Pyridine-d5	3.669	84	82826	7.855	ng/ul	0.00
7) Phenol-d5	6.987	99	56532	4.200	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.152	67	186465	23.371	ng/ul	0.00
11) 2-Chlorophenol-d4	7.346	132	176553	16.808	ng/ul	0.00
15) 4-Methylphenol-d8	8.528	113	103926	9.540	ng/ul	0.00
21) Nitrobenzene-d5	8.975	128	109906	24.939	ng/ul	0.00
24) 2-Nitrophenol-d4	9.693	143	109082	23.874	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.234	165	194285	20.795	ng/ul	0.00
31) 4-Chloroaniline-d4	10.751	131	259867	21.283	ng/ul	0.00
46) Dimethylphthalate-d6	13.863	166	682604	25.384	ng/ul	0.00
49) Acenaphthylene-d8	14.140	160	756342	25.543	ng/ul	0.00
54) 4-Nitrophenol-d4	14.681	143	15660m	3.548	ng/ul	0.02
60) Fluorene-d10	15.439	176	618599	27.405	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.569	200	97264	24.047	ng/ul	0.00
73) Anthracene-d10	17.292	188	1120710	31.602	ng/ul	0.00
81) Pyrene-d10	19.533	212	1432555	25.130	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.492	264	1722519	31.954	ng/ul	0.00

Target Compounds Qvalue

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

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