

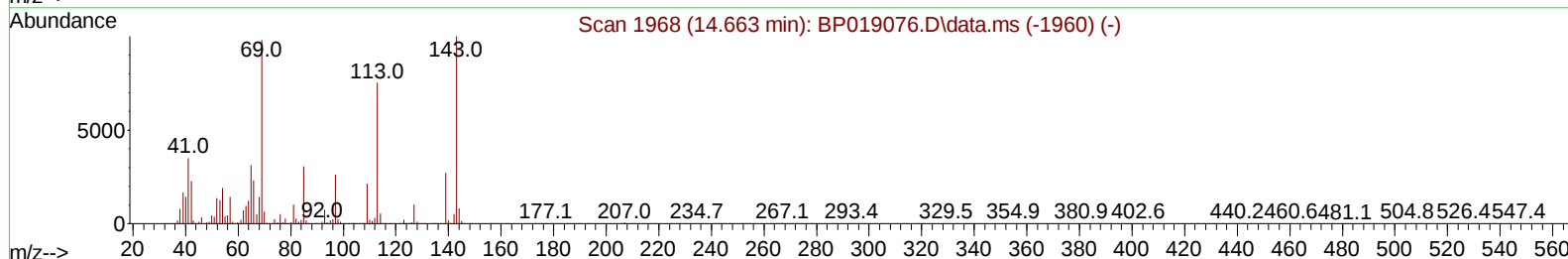
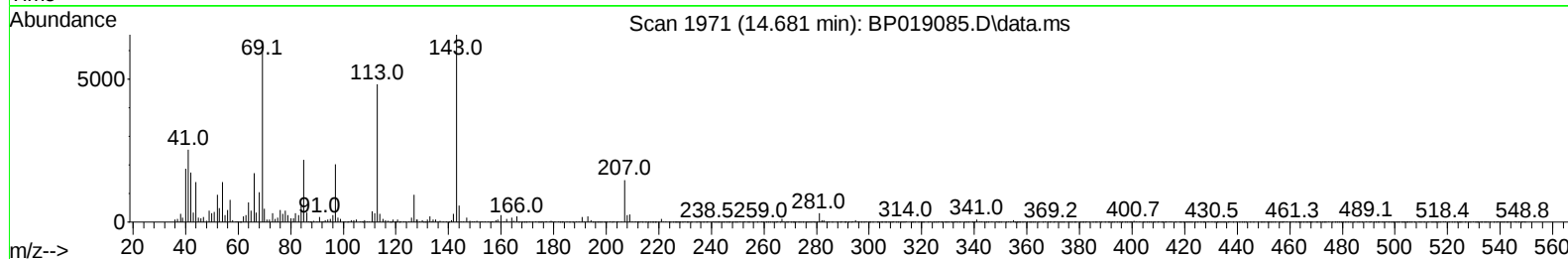
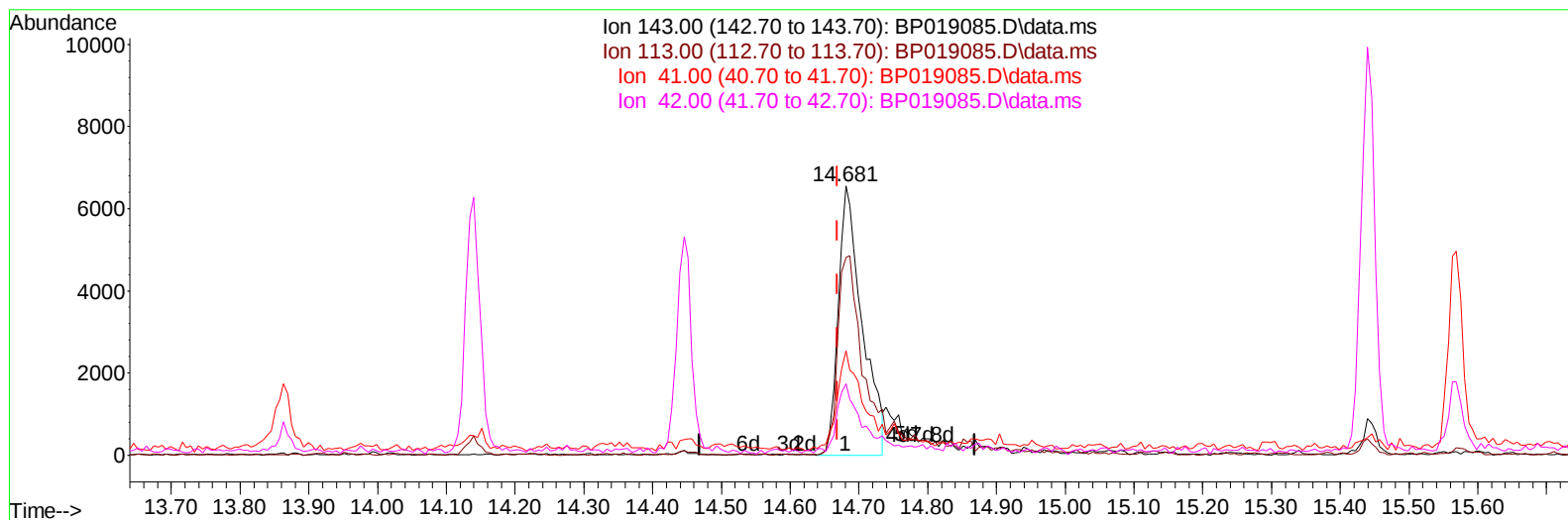
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122623\
 Data File : BP019085.D
 Acq On : 26 Dec 2023 16:04
 Operator : MA/JU
 Sample : 05976-02
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BH7T9

Manual IntegrationsAPPROVED

Quant Time: Dec 26 21:48:38 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120123.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Dec 21 04:46:12 2023
 Response via : Initial Calibration

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



TIC: BP019085.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.681min (+ 0.012) 3.69 ng/u1

response 15751

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	70.80	73.53
41.00	34.60	38.75
42.00	23.00	26.55

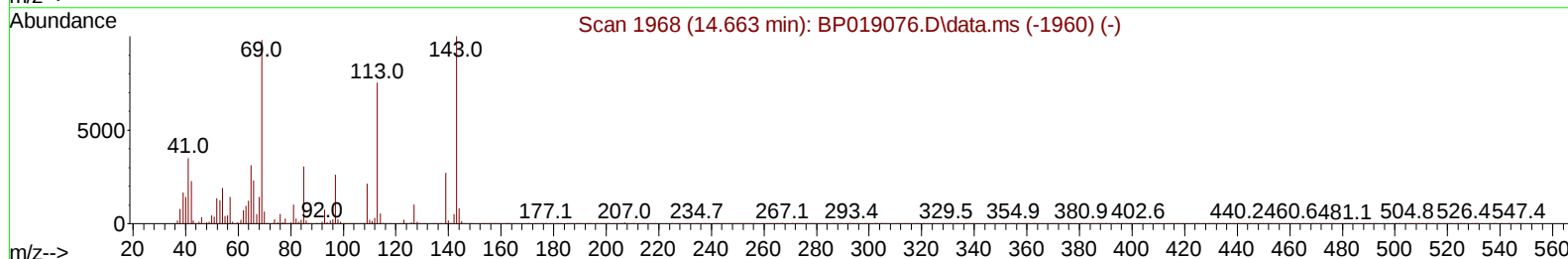
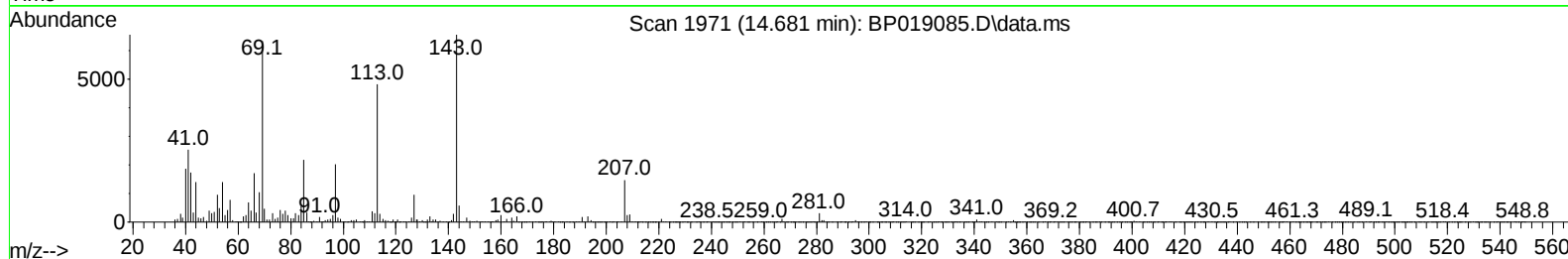
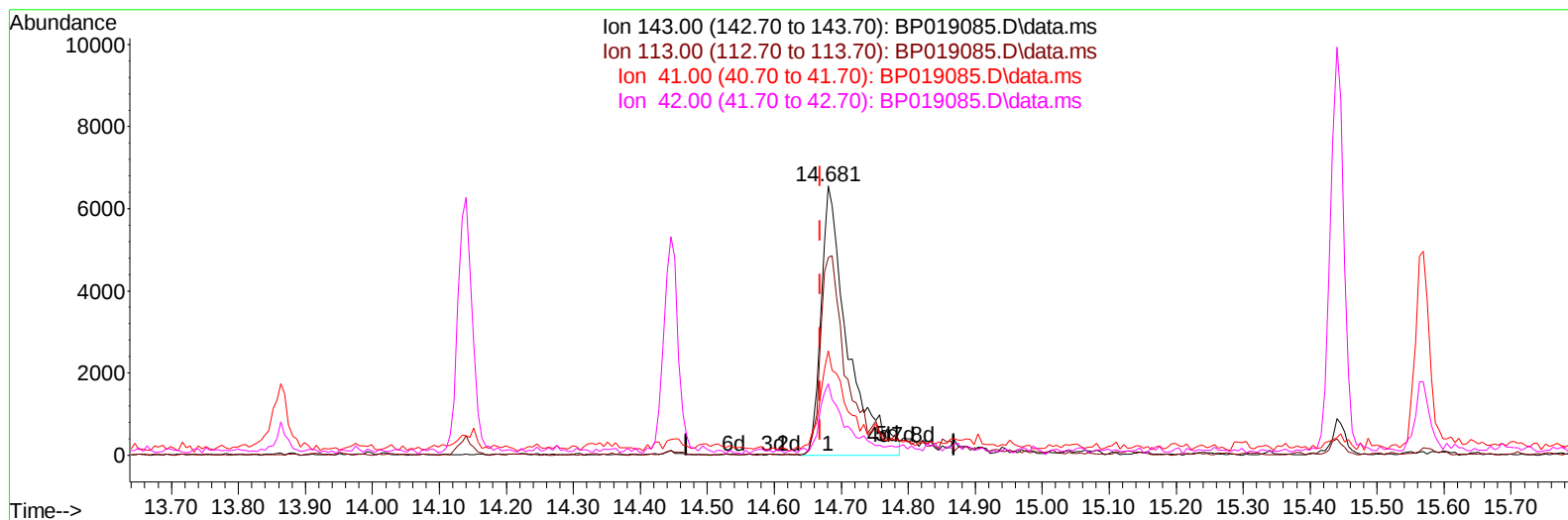
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TIC: BP019085.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.681min (+ 0.012) 4.21 ng/ul m

response 17957

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	70.80	73.53
41.00	34.60	38.75
42.00	23.00	26.55

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Manual IntegrationsAPPROVED

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

Quant Time: Dec 26 22:14:43 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120123.MA.M
 Quant Title : SVOA CALIBRATION
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.811	152	131314	20.000	ng/ul	-0.01
20) Naphthalene-d8	10.605	136	488204	20.000	ng/ul	-0.01
38) Acenaphthene-d10	14.446	164	291008	20.000	ng/ul	-0.01
64) Phenanthrene-d10	17.198	188	645791	20.000	ng/ul	-0.01
79) Chrysene-d12	21.286	240	688753	20.000	ng/ul	-0.01
88) Perylene-d12	23.645	264	863320	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.246	96	14982	3.888	ng/uL	0.00
4) Pyridine-d5	3.670	84	71257	6.904	ng/ul	0.00
7) Phenol-d5	6.993	99	61951	4.701	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.152	67	214150	27.419	ng/ul	-0.01
11) 2-Chlorophenol-d4	7.346	132	204340	19.872	ng/ul	0.00
15) 4-Methylphenol-d8	8.528	113	119972	11.250	ng/ul	-0.01
21) Nitrobenzene-d5	8.975	128	129294	29.756	ng/ul	-0.01
24) 2-Nitrophenol-d4	9.693	143	132244	29.355	ng/ul	-0.01
28) 2,4-Dichlorophenol-d3	10.234	165	231268	25.106	ng/ul	-0.01
31) 4-Chloroaniline-d4	10.752	131	297716	24.730	ng/ul	-0.01
46) Dimethylphthalate-d6	13.863	166	820393	31.587	ng/ul	-0.01
49) Acenaphthylene-d8	14.140	160	899647	31.457	ng/ul	-0.01
54) 4-Nitrophenol-d4	14.681	143	17957m	4.212	ng/ul	0.01
60) Fluorene-d10	15.440	176	732370	33.593	ng/ul	-0.01
65) 4,6-Dinitro-2-methylph...	15.569	200	116950	30.203	ng/ul	0.00
73) Anthracene-d10	17.298	188	1220419	35.947	ng/ul	0.00
81) Pyrene-d10	19.534	212	1567893	29.244	ng/ul	-0.01
92) Benzo(a)pyrene-d12	23.492	264	1872081	37.190	ng/ul	-0.01

Target Compounds Qvalue

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

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