

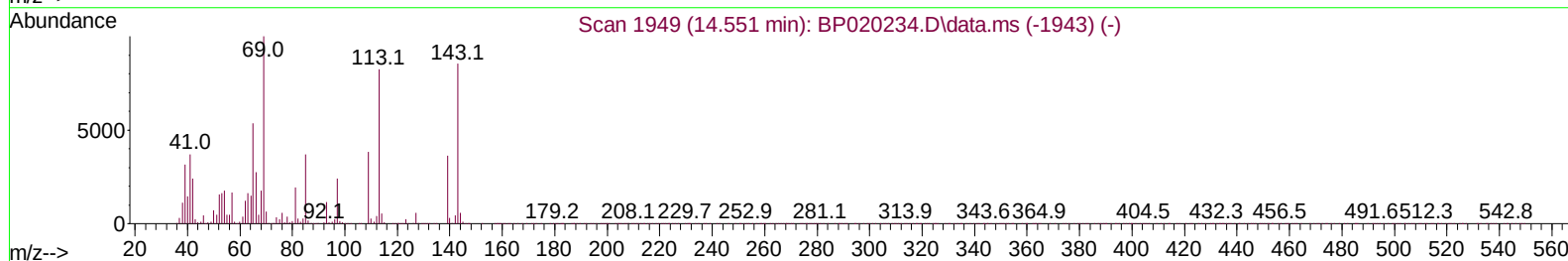
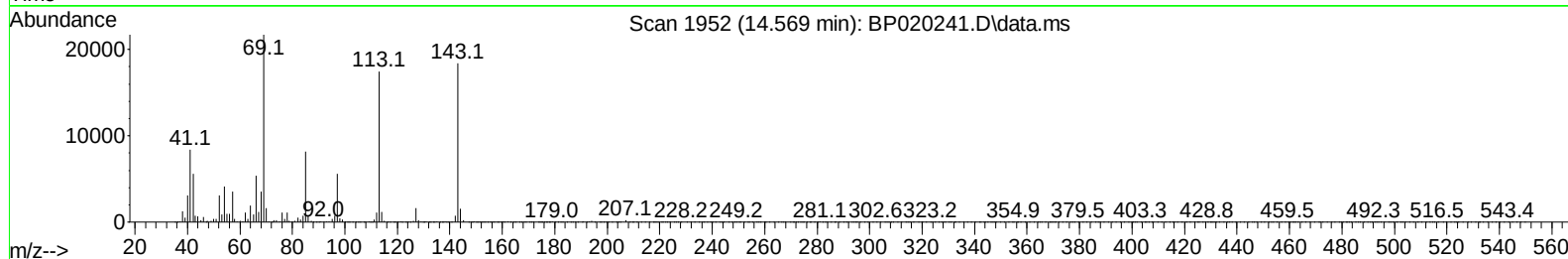
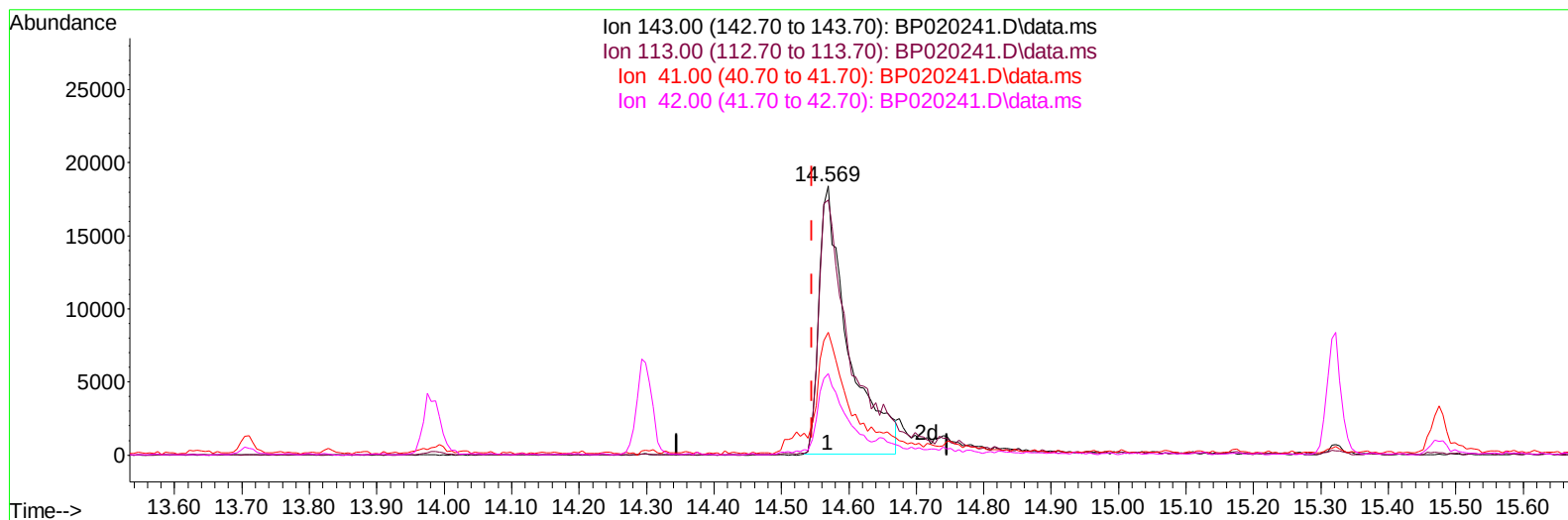
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP050824\
Data File : BP020241.D
Acq On : 08 May 2024 16:05
Operator : MA/JU
Sample : P2369-05
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A43M8

Manual IntegrationsAPPROVED

Quant Time: May 08 16:32:52 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 05/09/2024
Supervised By :mohammad ahmed 05/11/2024



TIC: BP020241.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.569min (+ 0.024) 16.63 ng/ul

response 55104

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	101.40	94.83
41.00	50.30	45.70
42.00	32.50	30.42

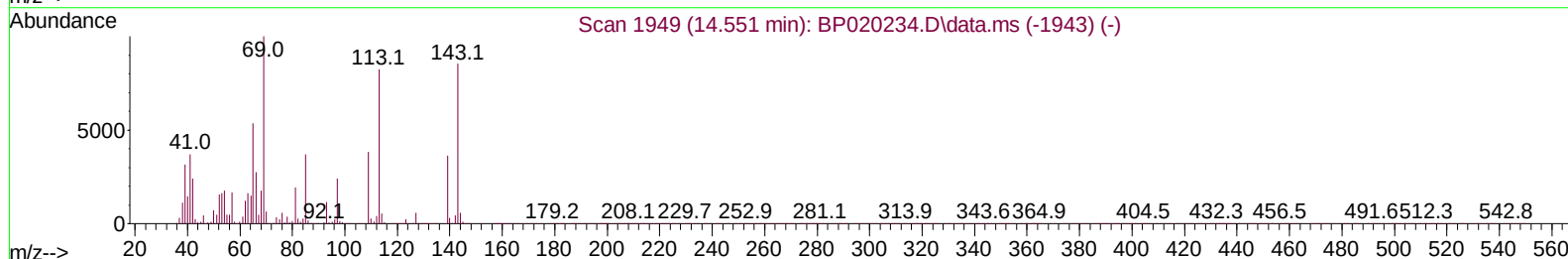
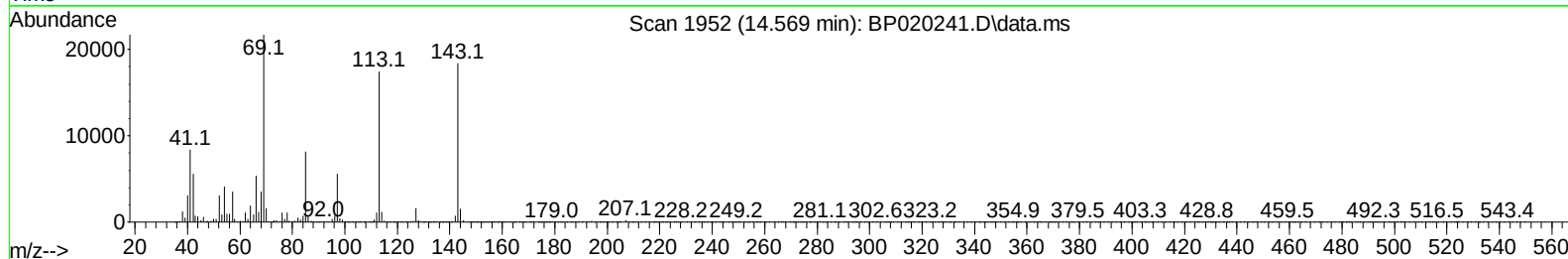
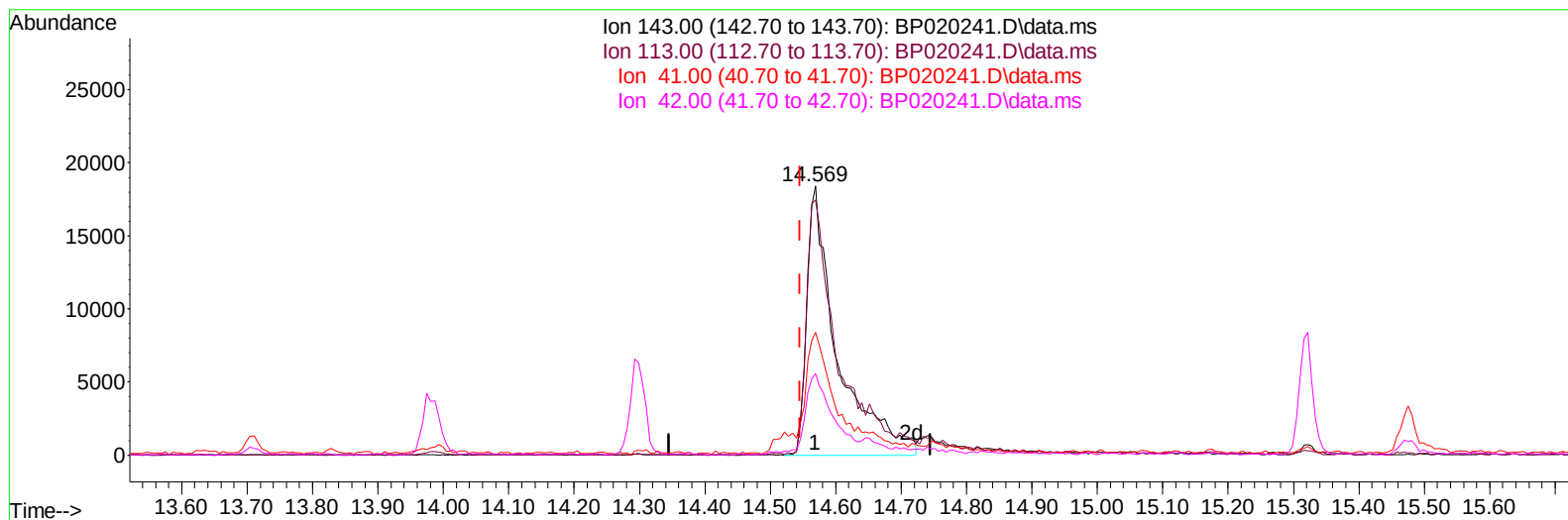
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TIC: BP020241.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.569min (+ 0.024) 18.10 ng/ul m

response 59979

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	101.40	94.83
41.00	50.30	45.70
42.00	32.50	30.42

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.616	152	93590	20.000	ng/ul	0.02
20) Naphthalene-d8	10.393	136	399677	20.000	ng/ul	0.02
38) Acenaphthene-d10	14.298	164	295280	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.134	188	670206	20.000	ng/ul	0.00
79) Chrysene-d12	21.639	240	576496	20.000	ng/ul	0.00
88) Perylene-d12	25.010	264	523378	20.000	ng/ul	0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.211	96	7502	3.385	ng/uL	0.00
4) Pyridine-d5	3.617	84	107369	16.675	ng/ul	0.01
7) Phenol-d5	6.811	99	148803	18.514	ng/ul	0.03
9) Bis-(2-Chloroethyl)eth...	6.969	67	95432	19.546	ng/ul	0.02
11) 2-Chlorophenol-d4	7.158	132	117713	20.487	ng/ul	0.02
15) 4-Methylphenol-d8	8.322	113	116896	17.989	ng/ul	0.01
21) Nitrobenzene-d5	8.775	128	61854	20.667	ng/ul	0.02
24) 2-Nitrophenol-d4	9.487	143	73007	21.301	ng/ul	0.02
28) 2,4-Dichlorophenol-d3	10.028	165	119352	19.083	ng/ul	0.02
31) 4-Chloroaniline-d4	10.552	131	176696	20.186	ng/ul	0.02
46) Dimethylphthalate-d6	13.710	166	486572	22.565	ng/ul	0.00
49) Acenaphthylene-d8	13.981	160	512468	21.514	ng/ul	0.00
54) 4-Nitrophenol-d4	14.569	143	59979m	18.098	ng/ul	0.02
60) Fluorene-d10	15.322	176	411465	22.878	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.475	200	66692	15.681	ng/ul	0.00
73) Anthracene-d10	17.240	188	648976	22.383	ng/ul	0.00
81) Pyrene-d10	19.651	212	800286	23.748	ng/ul	0.01
92) Benzo(a)pyrene-d12	24.780	264	581904	22.995	ng/ul	0.02
Target Compounds						
71) Pentachlorophenol	16.792	266	7784	1.532	ng/ul	Qvalue 93

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

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