

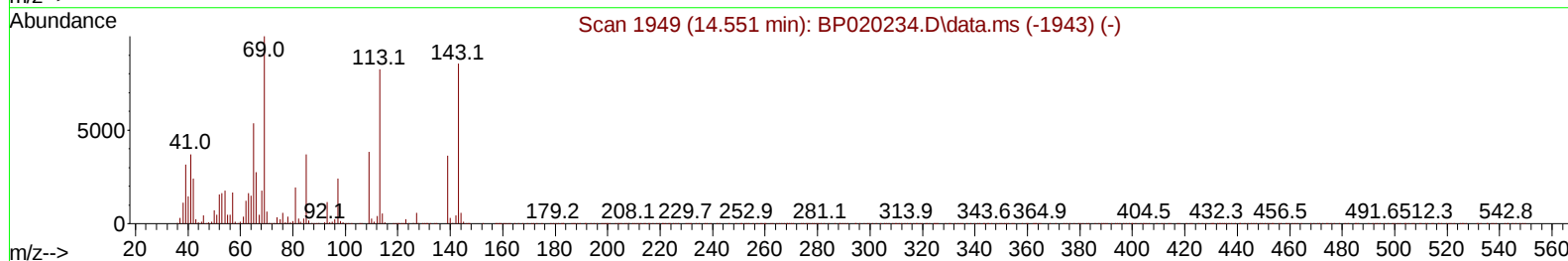
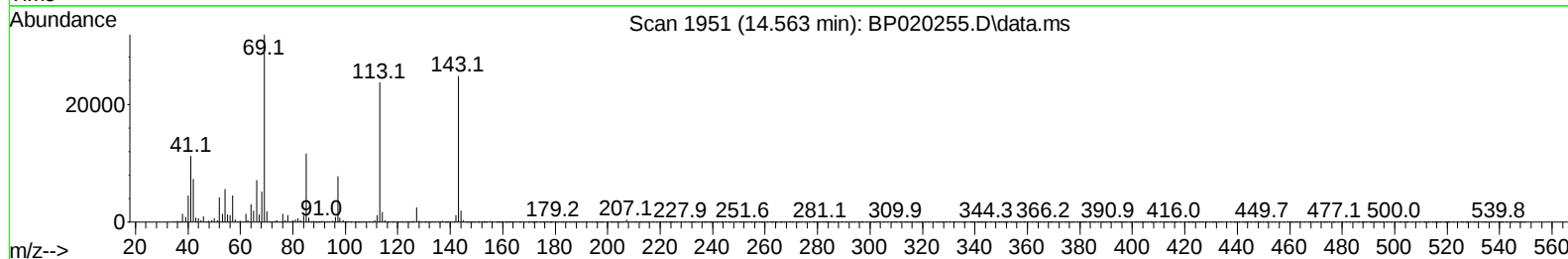
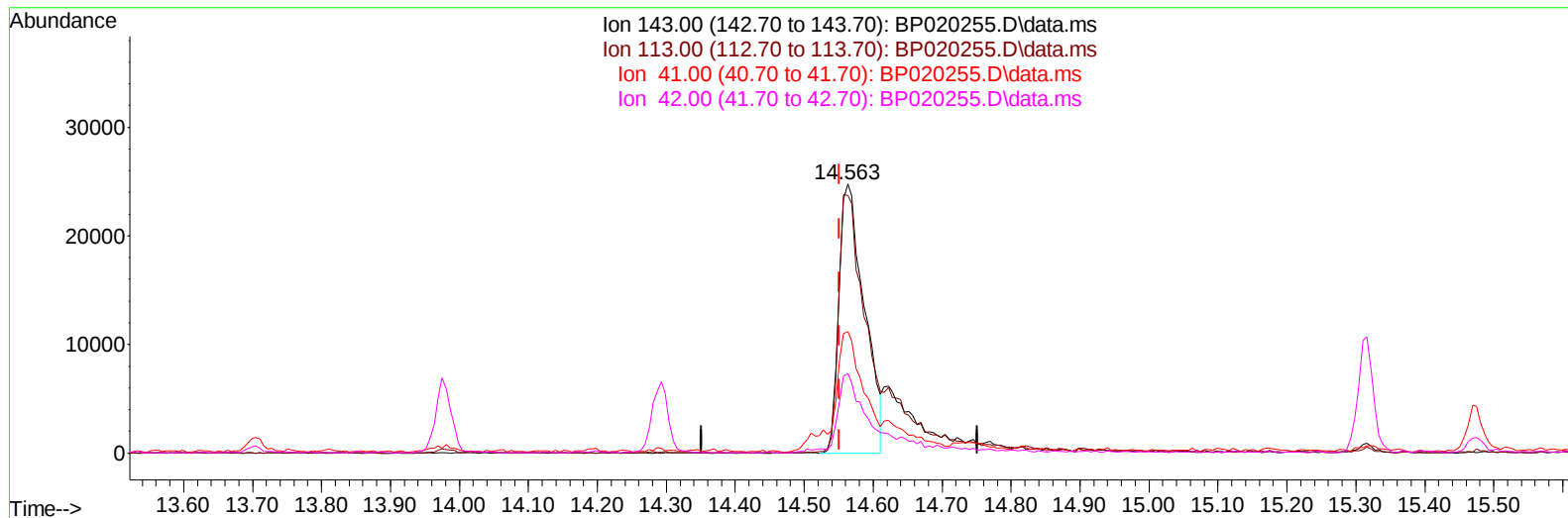
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP050824\
Data File : BP020255.D
Acq On : 09 May 2024 03:25
Operator : MA/JU
Sample : P2369-21
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A43P6

Manual IntegrationsAPPROVED

Quant Time: May 09 03:58:29 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP050124.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed May 08 22:57:55 2024
Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 05/09/2024
Supervised By :mohammad ahmed 05/11/2024



TIC: BP020255.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.563min (+ 0.012) 17.31 ng/ul

response 63313

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	101.40	95.80
41.00	50.30	45.17
42.00	32.50	29.50

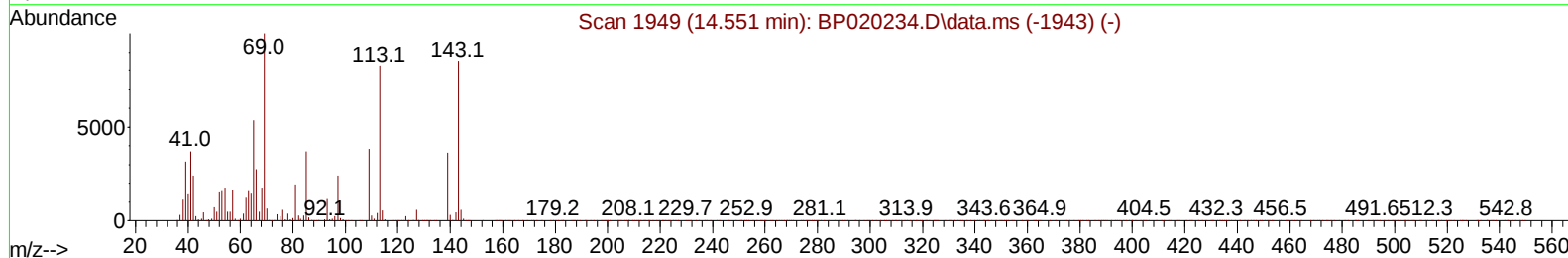
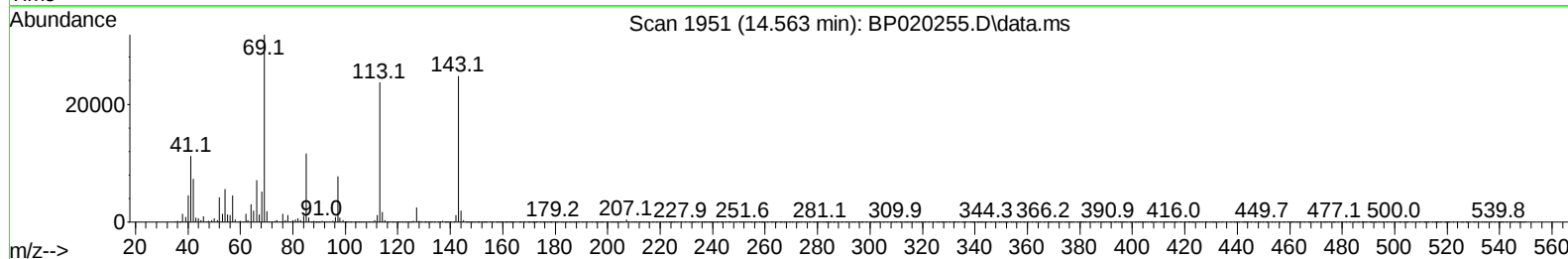
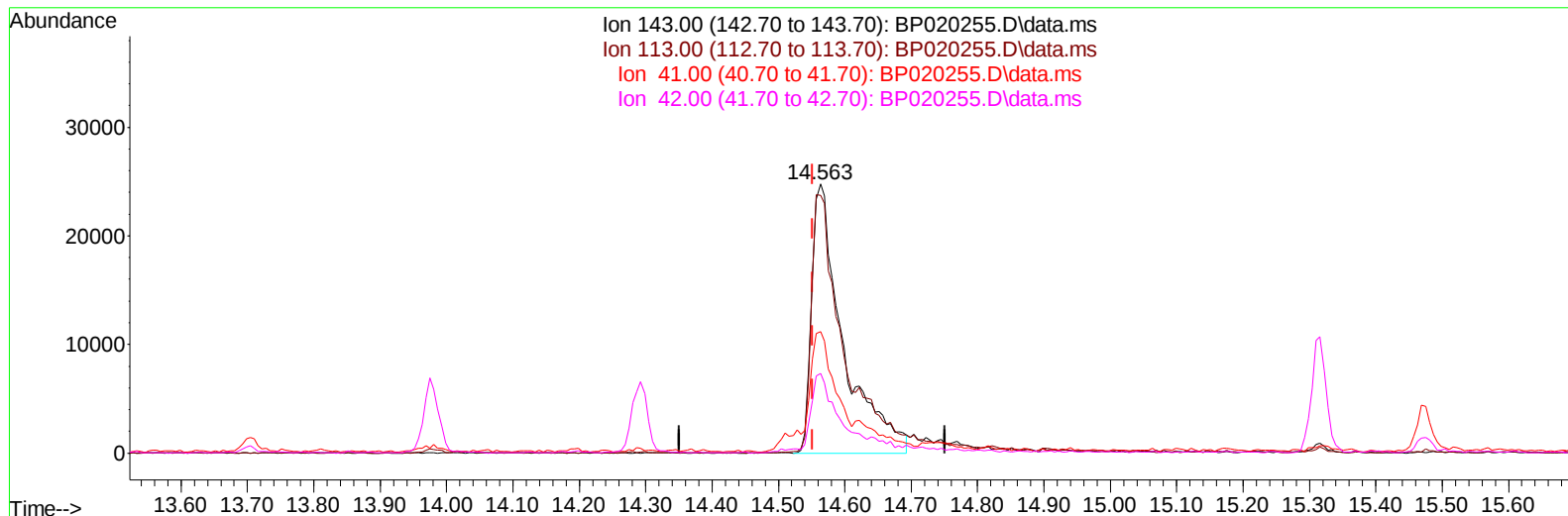
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(54) 4-Nitrophenol-d4 (S)

14.563min (+ 0.012) 22.24 ng/ul m

response 81318

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	101.40	95.80
41.00	50.30	45.17
42.00	32.50	29.50

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Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 05/09/2024
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Quant Time: May 09 03:59:12 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP050124.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed May 08 22:57:55 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.616	152	108076	20.000	ng/ul	0.00
20) Naphthalene-d8	10.387	136	461467	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.293	164	325841	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.134	188	654088	20.000	ng/ul	0.00
79) Chrysene-d12	21.628	240	463156	20.000	ng/ul	-0.01
88) Perylene-d12	24.992	264	521102	20.000	ng/ul	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.217	96	9538	3.727	ng/uL	0.00
4) Pyridine-d5	3.623	84	139862	18.810	ng/ul	0.00
7) Phenol-d5	6.805	99	202578	21.826	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.969	67	127630	22.637	ng/ul	0.00
11) 2-Chlorophenol-d4	7.158	132	158179	23.840	ng/ul	0.00
15) 4-Methylphenol-d8	8.328	113	179510	23.922	ng/ul	0.00
21) Nitrobenzene-d5	8.781	128	87382	25.287	ng/ul	0.00
24) 2-Nitrophenol-d4	9.493	143	101263	25.590	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.028	165	186138	25.776	ng/ul	0.00
31) 4-Chloroaniline-d4	10.552	131	252584	24.992	ng/ul	0.00
46) Dimethylphthalate-d6	13.704	166	713156	29.971	ng/ul	0.00
49) Acenaphthylene-d8	13.981	160	746503	28.400	ng/ul	0.00
54) 4-Nitrophenol-d4	14.563	143	81318m	22.236	ng/ul	0.01
60) Fluorene-d10	15.316	176	581802	29.314	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.475	200	87970	21.194	ng/ul	0.00
73) Anthracene-d10	17.240	188	841348	29.734	ng/ul	0.00
81) Pyrene-d10	19.645	212	873037	32.247	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.763	264	743519	29.510	ng/ul	-0.01

Target Compounds Qvalue

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

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