

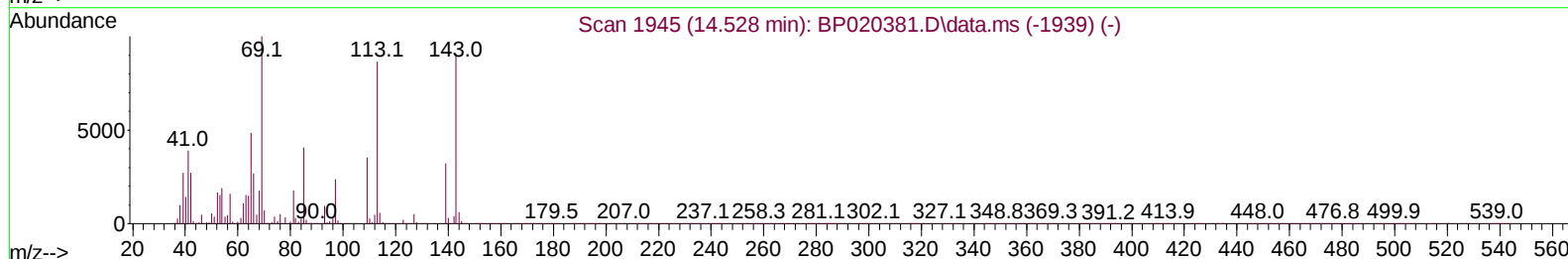
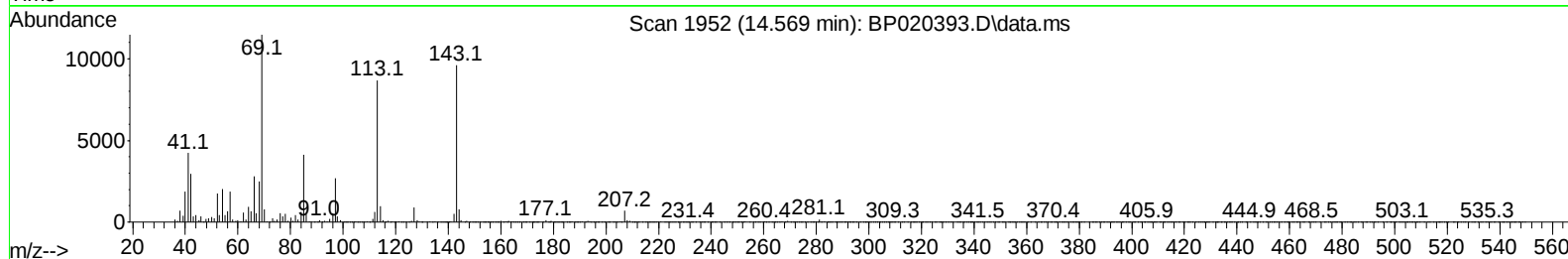
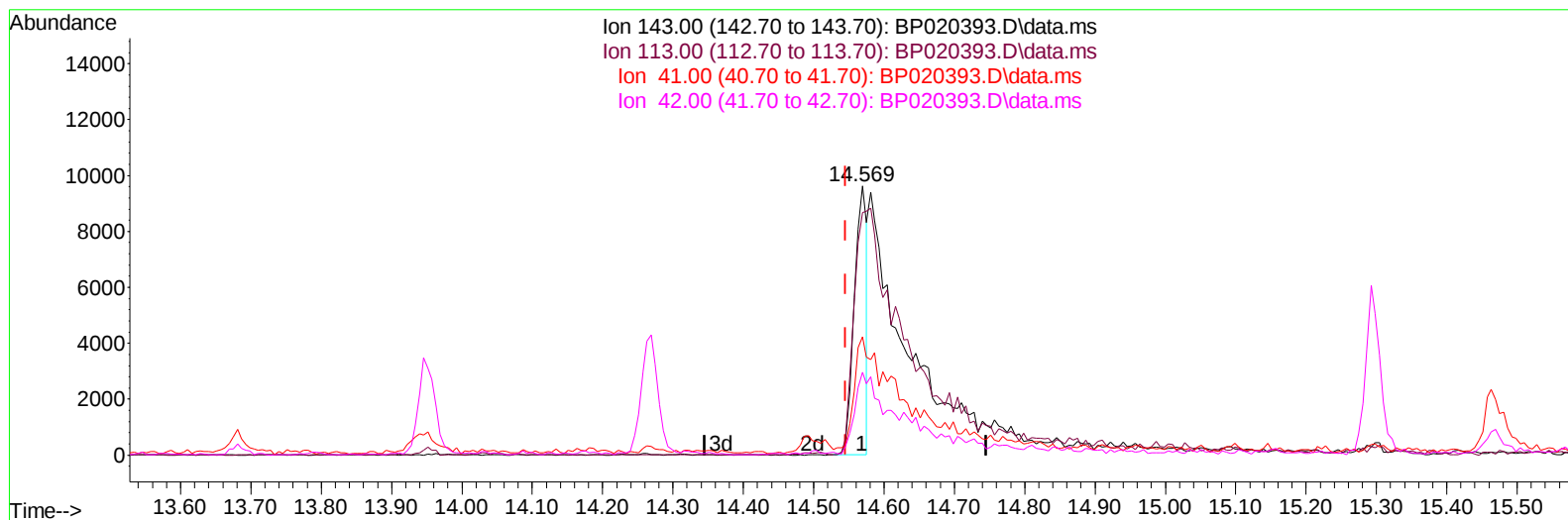
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051424\
Data File : BP020393.D
Acq On : 15 May 2024 09:40
Operator : MA/JU
Sample : P2372-21
Misc :
ALS Vial : 27 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A43X1

Manual IntegrationsAPPROVED

Quant Time: May 15 10:39:53 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP050124.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue May 14 23:58:44 2024
Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 05/15/2024
Supervised By :mohammad ahmed 05/17/2024



TIC: BP020393.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.569min (+ 0.024) 5.13 ng/u1

response 12158

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	101.40	90.27
41.00	50.30	44.03
42.00	32.50	30.82

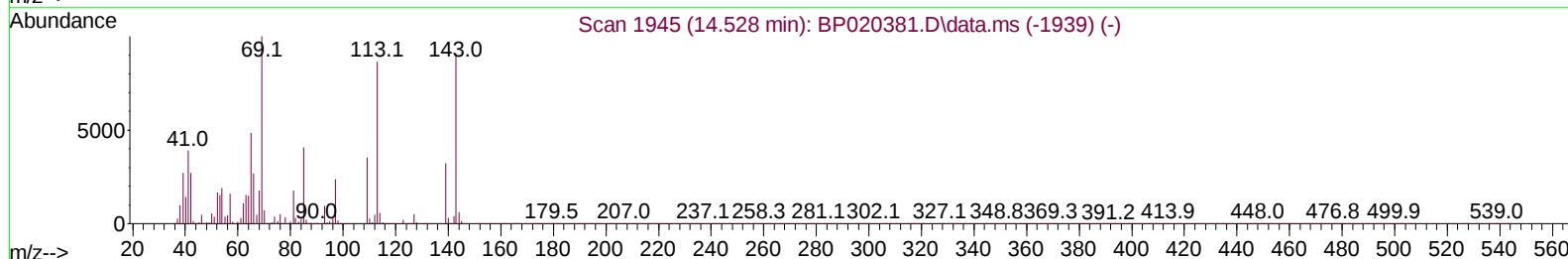
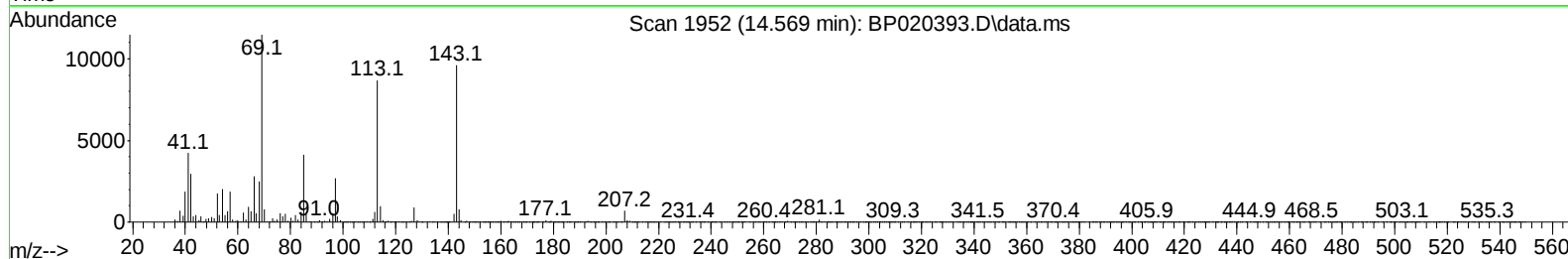
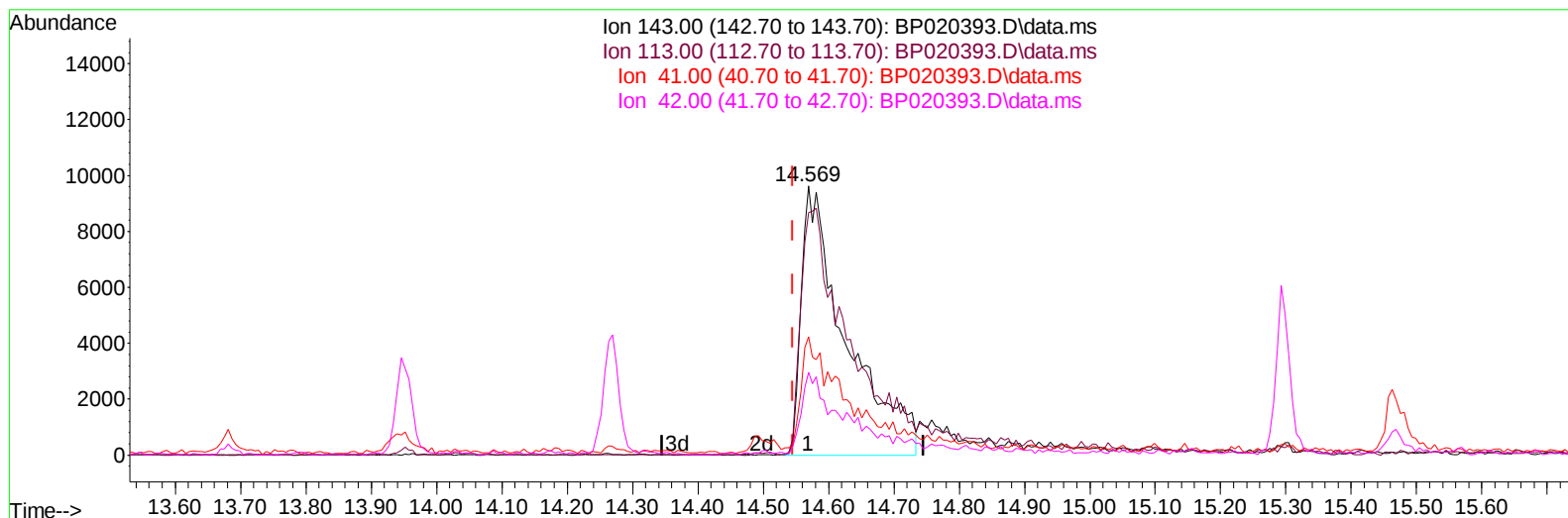
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TIC: BP020393.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.569min (+ 0.024) 19.22 ng/ul m

response 45544

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	101.40	90.27
41.00	50.30	44.03
42.00	32.50	30.82

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Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 05/15/2024
 Supervised By :mohammad ahmed 05/17/2024

Quant Time: May 15 10:42:07 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP050124.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue May 14 23:58:44 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.593	152	73309	20.000	ng/ul	0.00
20) Naphthalene-d8	10.357	136	296052	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.269	164	211086	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.116	188	506850	20.000	ng/ul	0.00
79) Chrysene-d12	21.610	240	552016	20.000	ng/ul	0.00
88) Perylene-d12	24.951	264	630266	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.187	96	7190	4.142	ng/uL	0.00
4) Pyridine-d5	3.593	84	87864	17.421	ng/ul	0.00
7) Phenol-d5	6.787	99	118173	18.770	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.946	67	72825	19.042	ng/ul	0.00
11) 2-Chlorophenol-d4	7.134	132	95866	21.300	ng/ul	0.00
15) 4-Methylphenol-d8	8.305	113	94215	18.510	ng/ul	0.00
21) Nitrobenzene-d5	8.752	128	47016	21.208	ng/ul	0.00
24) 2-Nitrophenol-d4	9.463	143	54173	21.339	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.010	165	95328	20.577	ng/ul	0.00
31) 4-Chloroaniline-d4	10.534	131	126363	19.489	ng/ul	0.00
46) Dimethylphthalate-d6	13.681	166	370003	24.003	ng/ul	0.00
49) Acenaphthylene-d8	13.951	160	388185	22.797	ng/ul	0.00
54) 4-Nitrophenol-d4	14.569	143	45544m	19.224	ng/ul	0.02
60) Fluorene-d10	15.298	176	314789	24.483	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.463	200	51416	15.986	ng/ul	0.00
73) Anthracene-d10	17.228	188	544386	24.828	ng/ul	0.01
81) Pyrene-d10	19.628	212	696339	21.580	ng/ul	0.01
92) Benzo(a)pyrene-d12	24.739	264	722778	23.718	ng/ul	0.02

Target Compounds Qvalue

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

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