

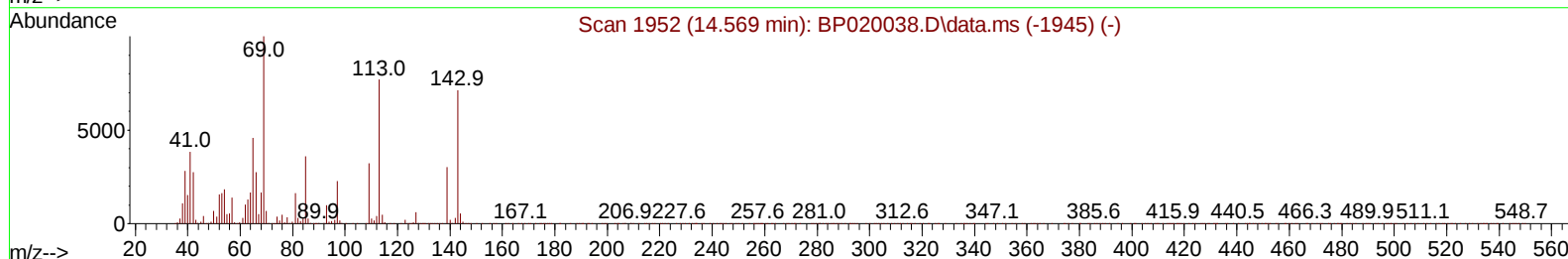
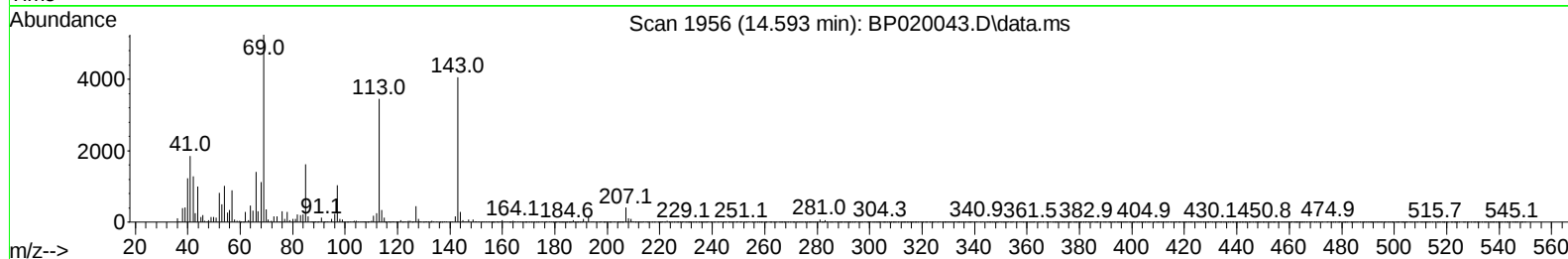
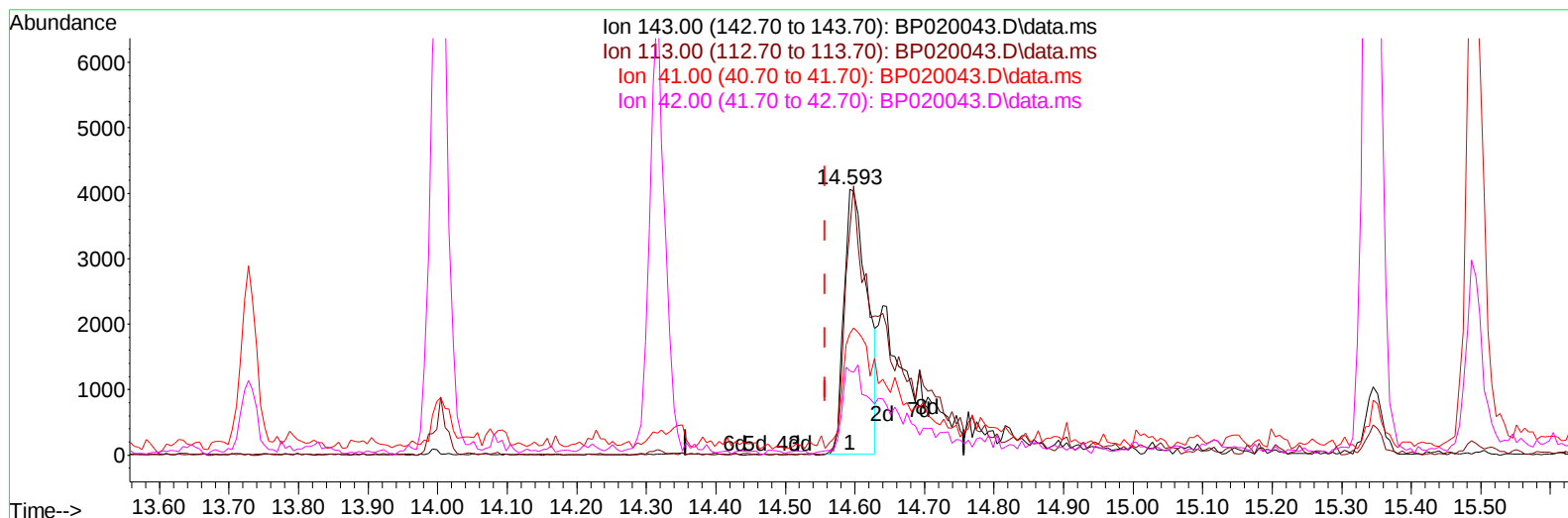
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042024\
Data File : BP020043.D
Acq On : 20 Apr 2024 14:21
Operator : MA/JU
Sample : P2161-06
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0RB3

Manual IntegrationsAPPROVED

Quant Time: Apr 21 23:42:49 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP041924.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Sat Apr 20 01:32:50 2024
Response via : Initial Calibration

Reviewed By :Yogesh Patel 04/24/2024
Supervised By :mohammad ahmed 04/30/2024



TIC: BP020043.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.593min (+ 0.035) 3.55 ng/u1

response 9322

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	97.80	84.96
41.00	50.70	45.56
42.00	34.00	31.55

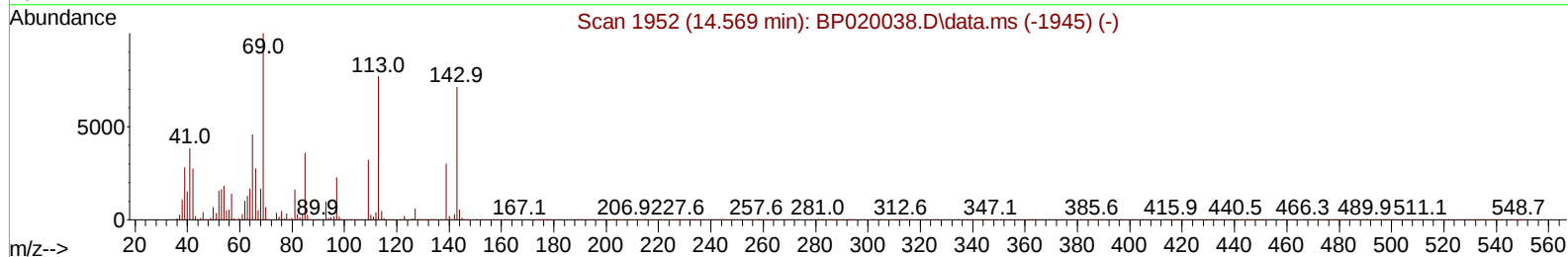
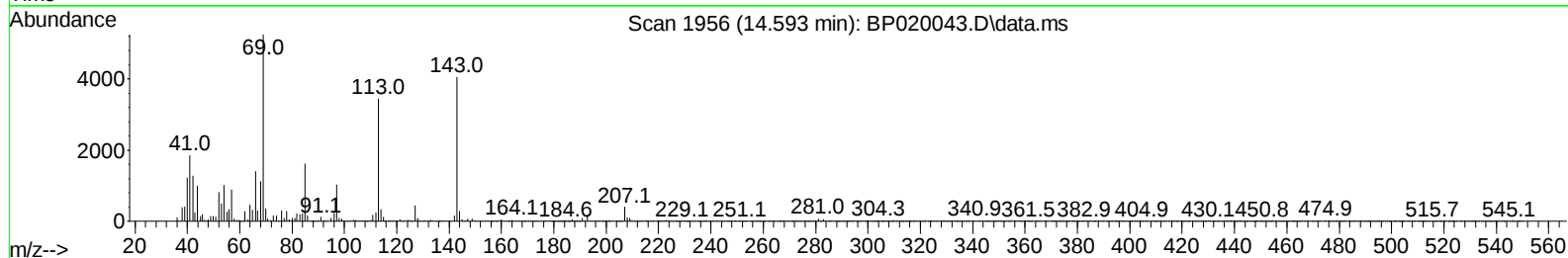
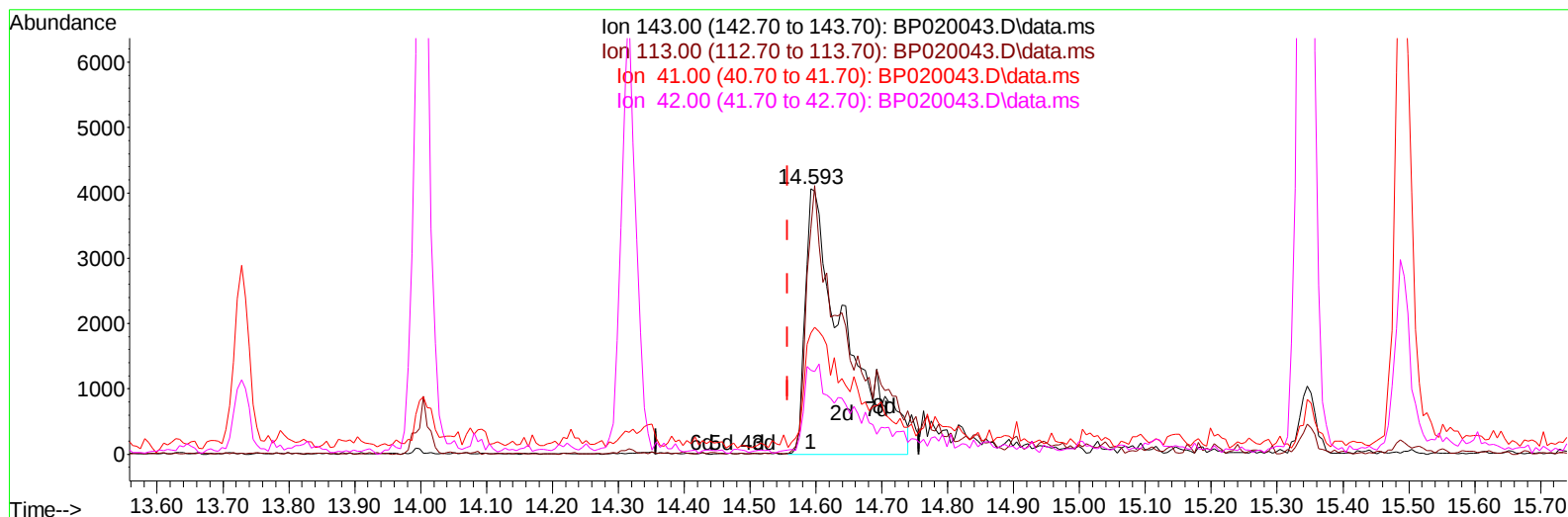
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TIC: BP020043.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.593min (+ 0.035) 6.55 ng/ul m

response 17198

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	97.80	84.96
41.00	50.70	45.56
42.00	34.00	31.55

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 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_P
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Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 04/24/2024
 Supervised By :mohammad ahmed 04/30/2024

Quant Time: Apr 21 23:58:12 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP041924.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Sat Apr 20 01:32:50 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.640	152	77213	20.000	ng/ul	0.00
20) Naphthalene-d8	10.416	136	348899	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.316	164	252321	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.157	188	525558	20.000	ng/ul	0.00
79) Chrysene-d12	21.669	240	397306	20.000	ng/ul	0.01
88) Perylene-d12	25.057	264	445282	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.234	96	9191	4.877	ng/uL	0.00
4) Pyridine-d5	3.640	84	36233	6.532	ng/ul	0.00
7) Phenol-d5	6.828	99	51752	7.690	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.993	67	214472	45.597	ng/ul	0.00
11) 2-Chlorophenol-d4	7.181	132	157798	34.343	ng/ul	0.00
15) 4-Methylphenol-d8	8.346	113	105946	20.155	ng/ul	0.00
21) Nitrobenzene-d5	8.799	128	116844	46.557	ng/ul	0.00
24) 2-Nitrophenol-d4	9.511	143	125008	46.080	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.046	165	197291	36.659	ng/ul	0.00
31) 4-Chloroaniline-d4	10.575	131	203922	28.384	ng/ul	0.00
46) Dimethylphthalate-d6	13.728	166	927295	53.775	ng/ul	0.00
49) Acenaphthylene-d8	14.004	160	955872	46.508	ng/ul	0.00
54) 4-Nitrophenol-d4	14.593	143	17198m	6.552	ng/ul	0.04
60) Fluorene-d10	15.346	176	784952	51.137	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.493	200	140504	45.761	ng/ul	0.00
73) Anthracene-d10	17.269	188	1233877	52.404	ng/ul	0.00
81) Pyrene-d10	19.669	212	1282634	60.072	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.821	264	1164801	53.549	ng/ul	0.01

Target Compounds Qvalue

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

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