

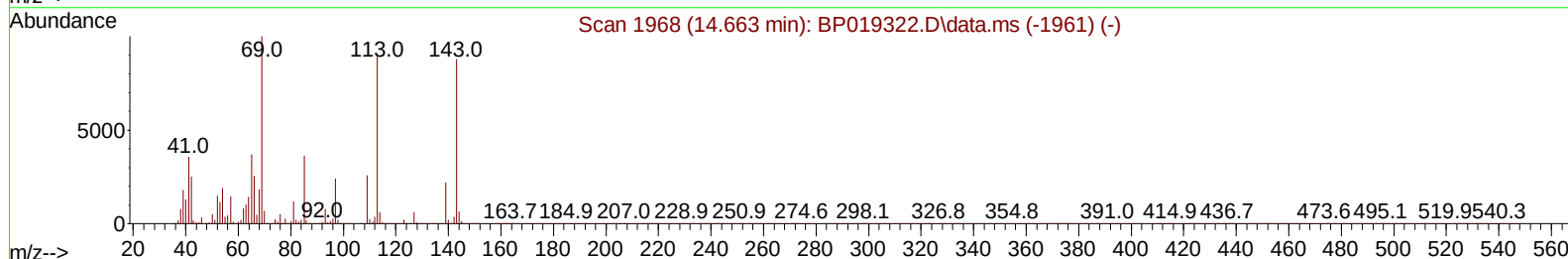
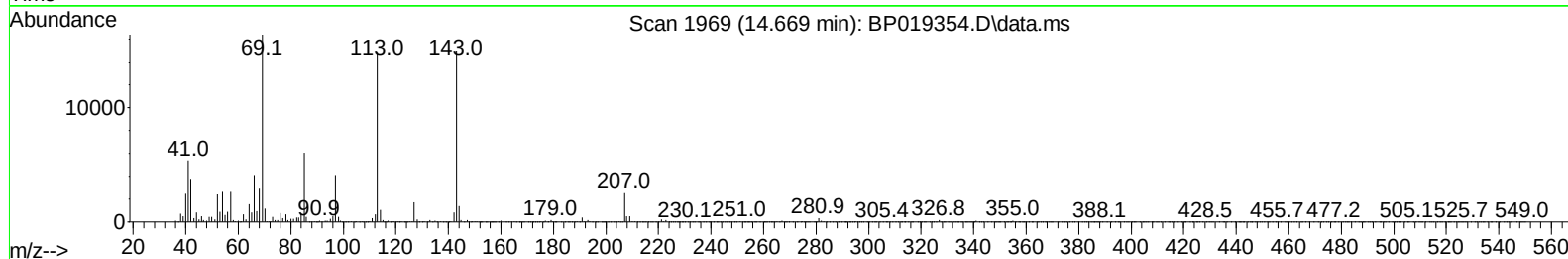
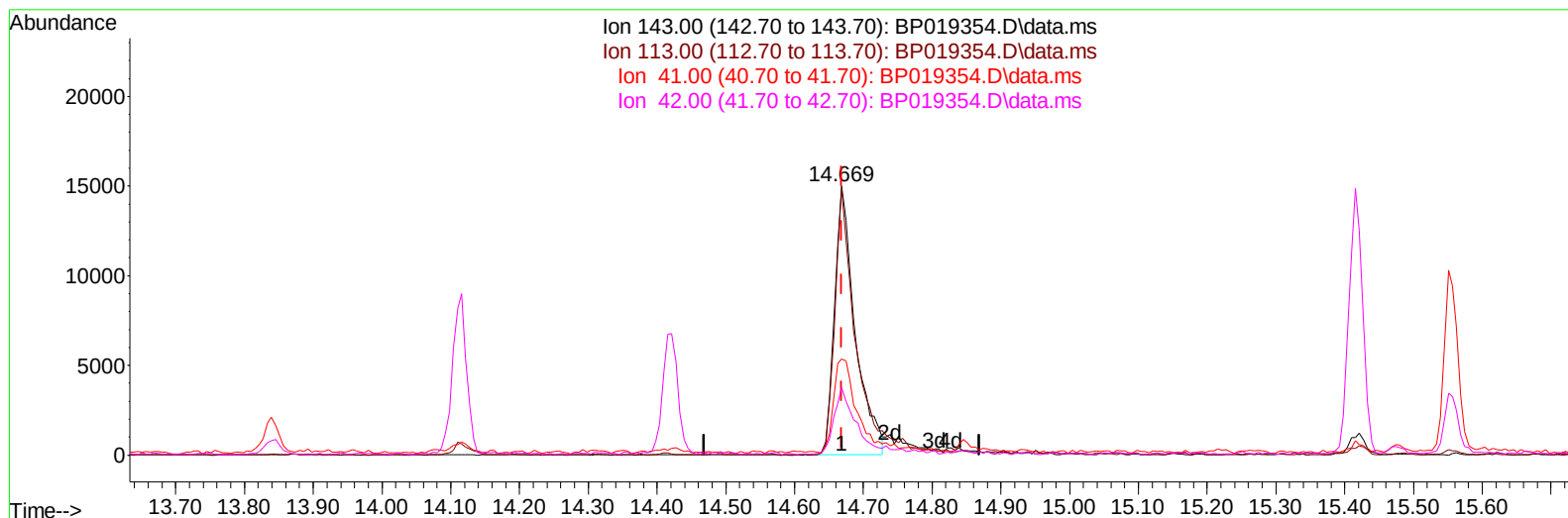
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP011224\
Data File : BP019354.D
Acq On : 13 Jan 2024 05:27
Operator : MA/JU
Sample : P1091-11
Misc :
ALS Vial : 34 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0AX7

Manual IntegrationsAPPROVED

Quant Time: Jan 13 05:54:42 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP010824.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue Jan 09 03:05:24 2024
Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 01/15/2024
Supervised By :mohammad ahmed 01/16/2024



TIC: BP019354.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.669min (-0.000) 6.64 ng/u1

response 31385

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	98.20	99.09
41.00	39.60	35.82
42.00	25.70	25.13

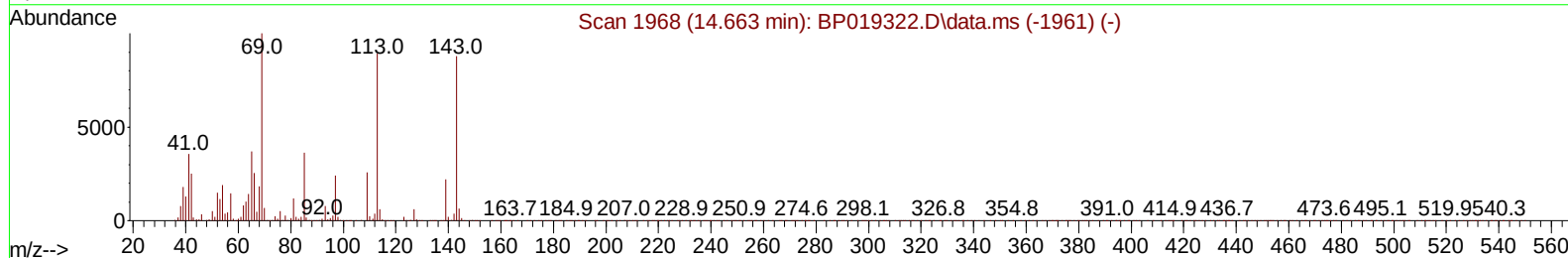
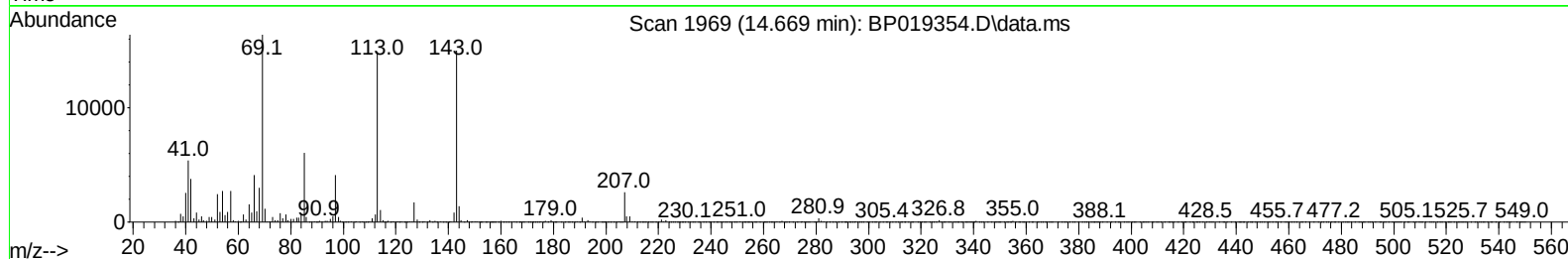
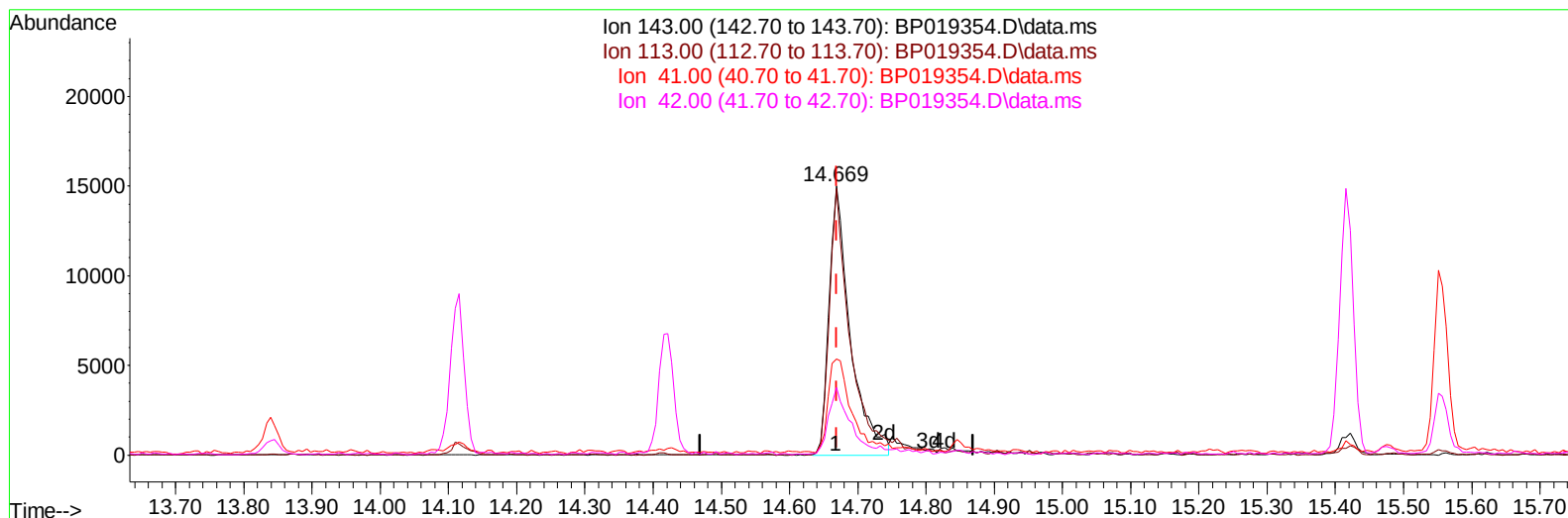
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TIC: BP019354.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.669min (-0.000) 6.86 ng/ul m

response 32409

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	98.20	99.09
41.00	39.60	35.82
42.00	25.70	25.13

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.775	152	110090	20.000	ng/ul	-0.01
20) Naphthalene-d8	10.569	136	458426	20.000	ng/ul	-0.01
38) Acenaphthene-d10	14.422	164	348171	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.180	188	841131	20.000	ng/ul	0.00
79) Chrysene-d12	21.286	240	882277	20.000	ng/ul	0.00
88) Perylene-d12	23.639	264	909051	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.216	96	12801	4.397	ng/uL	0.00
4) Pyridine-d5	3.640	84	80563	9.569	ng/ul	0.00
7) Phenol-d5	6.963	99	61769	5.870	ng/ul	-0.01
9) Bis-(2-Chloroethyl)eth...	7.116	67	209006	33.209	ng/ul	-0.01
11) 2-Chlorophenol-d4	7.310	132	167477	23.171	ng/ul	0.00
15) 4-Methylphenol-d8	8.504	113	75439	9.037	ng/ul	-0.01
21) Nitrobenzene-d5	8.945	128	127664	33.102	ng/ul	0.00
24) 2-Nitrophenol-d4	9.663	143	148519	33.143	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.204	165	240366	26.550	ng/ul	0.00
31) 4-Chloroaniline-d4	10.728	131	84565	7.395	ng/ul	0.00
46) Dimethylphthalate-d6	13.839	166	1064149	34.595	ng/ul	0.00
49) Acenaphthylene-d8	14.116	160	1079757	32.747	ng/ul	0.00
54) 4-Nitrophenol-d4	14.669	143	32409m	6.861	ng/ul	0.00
60) Fluorene-d10	15.416	176	921474	36.590	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.557	200	223006	35.911	ng/ul	0.00
73) Anthracene-d10	17.274	188	1390088	33.136	ng/ul	-0.01
81) Pyrene-d10	19.527	212	1951145	33.915	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.486	264	1939470	39.237	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.252	88	45871	15.211	ng/uL	97

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

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