

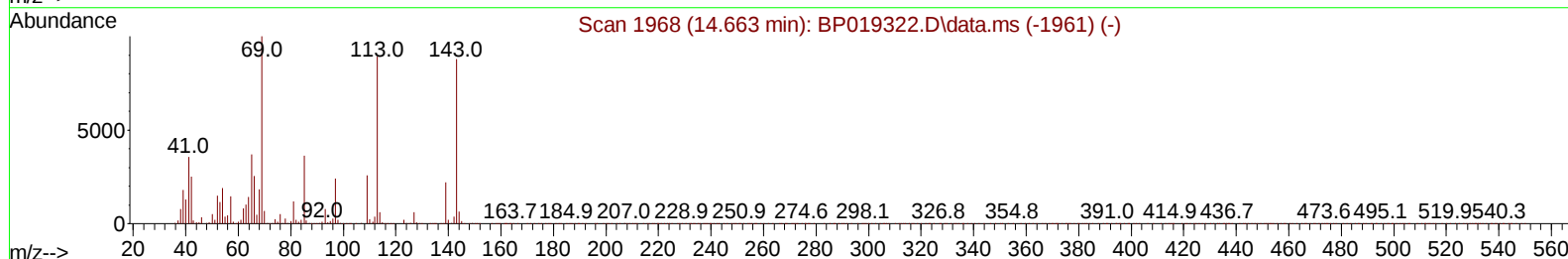
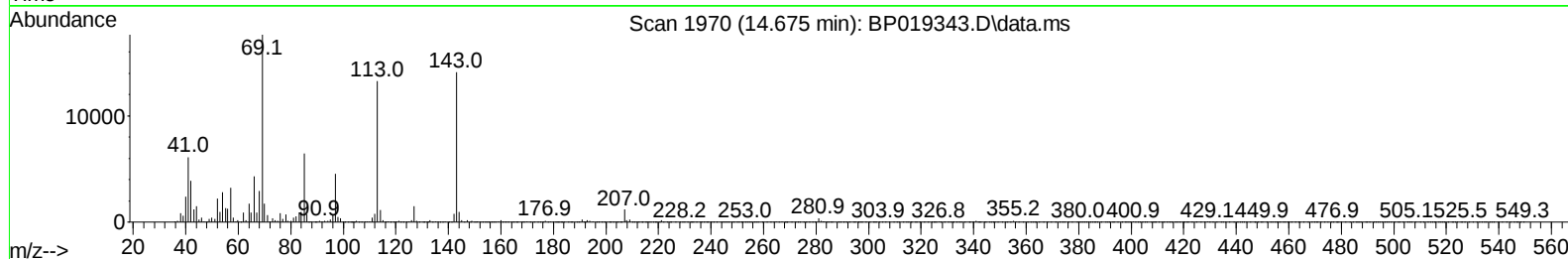
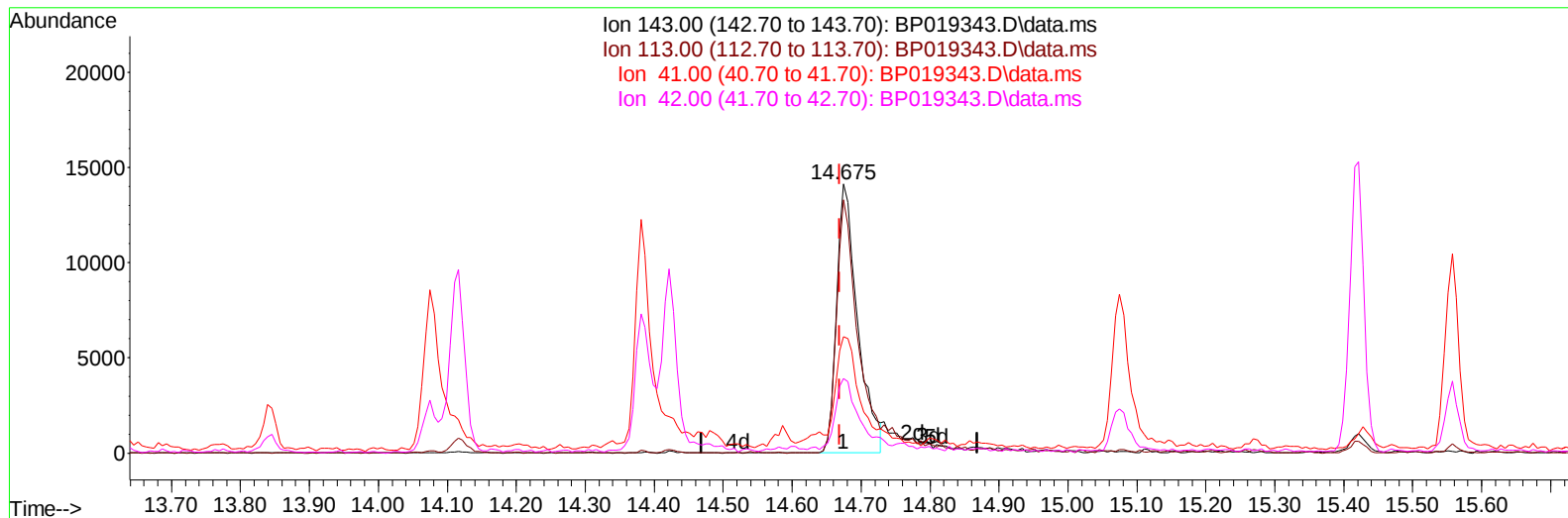
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP011224\
Data File : BP019343.D
Acq On : 12 Jan 2024 22:38
Operator : MA/JU
Sample : P1091-13
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0AY9

Manual IntegrationsAPPROVED

Quant Time: Jan 12 23:10:24 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: BP019343.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.675min (+ 0.006) 5.46 ng/u1

response 29970

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	98.20	94.13
41.00	39.60	43.37
42.00	25.70	27.69

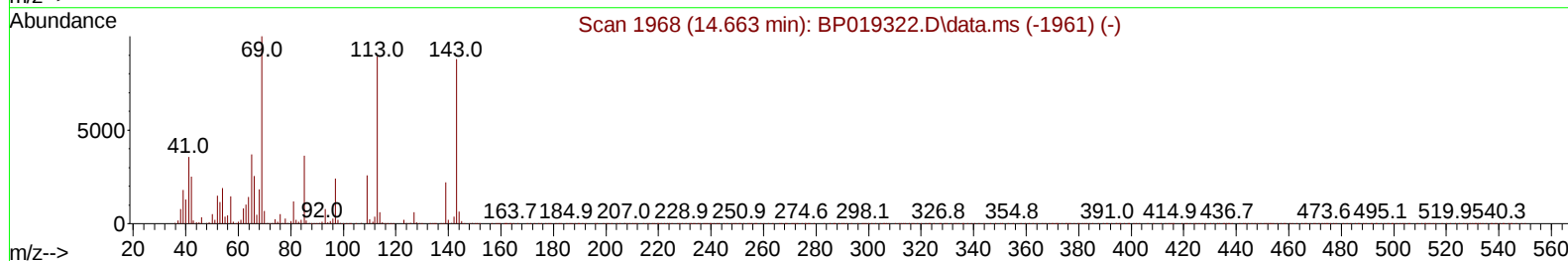
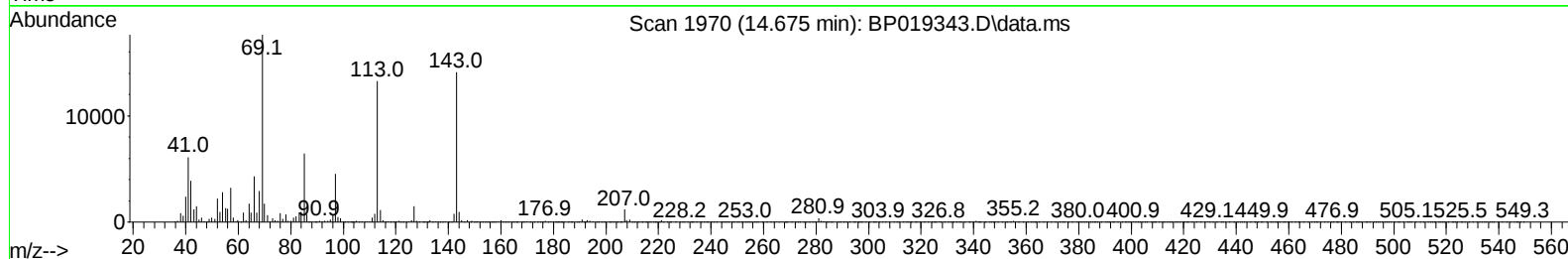
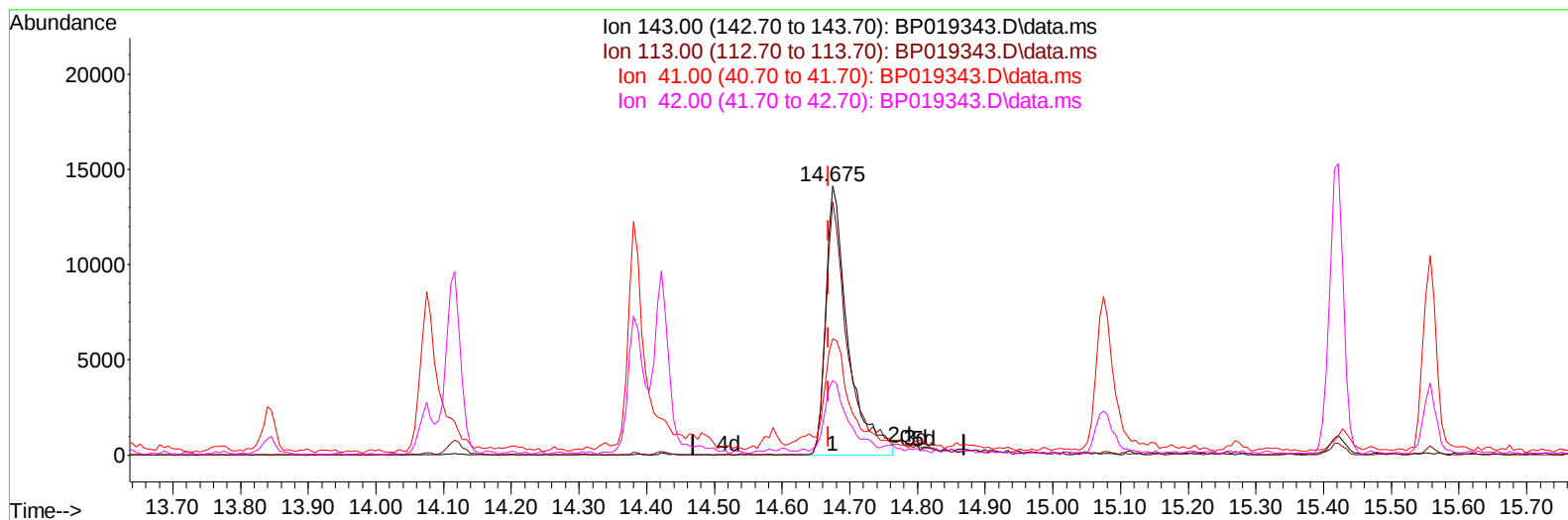
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TIC: BP019343.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.675min (+ 0.006) 5.88 ng/ul m

response 32264

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	98.20	94.13
41.00	39.60	43.37
42.00	25.70	27.69

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Quant Time: Jan 12 23:11:42 2024
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.775	152	134396	20.000	ng/ul	-0.01
20) Naphthalene-d8	10.575	136	550455	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.422	164	404634	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.181	188	1036802	20.000	ng/ul	0.00
79) Chrysene-d12	21.286	240	1048599	20.000	ng/ul	0.00
88) Perylene-d12	23.639	264	1059873	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.217	96	14992	4.218	ng/uL	0.00
4) Pyridine-d5	3.640	84	100385	9.767	ng/ul	0.00
7) Phenol-d5	6.969	99	81044	6.309	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.122	67	229700	29.896	ng/ul	0.00
11) 2-Chlorophenol-d4	7.316	132	206407	23.392	ng/ul	0.00
15) 4-Methylphenol-d8	8.505	113	153008	15.015	ng/ul	-0.01
21) Nitrobenzene-d5	8.946	128	137402	29.670	ng/ul	0.00
24) 2-Nitrophenol-d4	9.663	143	161099	29.940	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.204	165	291176	26.785	ng/ul	0.00
31) 4-Chloroaniline-d4	10.728	131	328179	23.899	ng/ul	0.00
46) Dimethylphthalate-d6	13.845	166	1112046	31.108	ng/ul	0.00
49) Acenaphthylene-d8	14.116	160	1157477	30.205	ng/ul	0.00
54) 4-Nitrophenol-d4	14.675	143	32264m	5.877	ng/ul	0.00
60) Fluorene-d10	15.422	176	972468	33.227	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.557	200	231440	30.236	ng/ul	0.00
73) Anthracene-d10	17.281	188	1728733	33.431	ng/ul	0.00
81) Pyrene-d10	19.527	212	2232282	32.647	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.486	264	2010741	34.890	ng/ul	0.00

Target Compounds Qvalue

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

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