

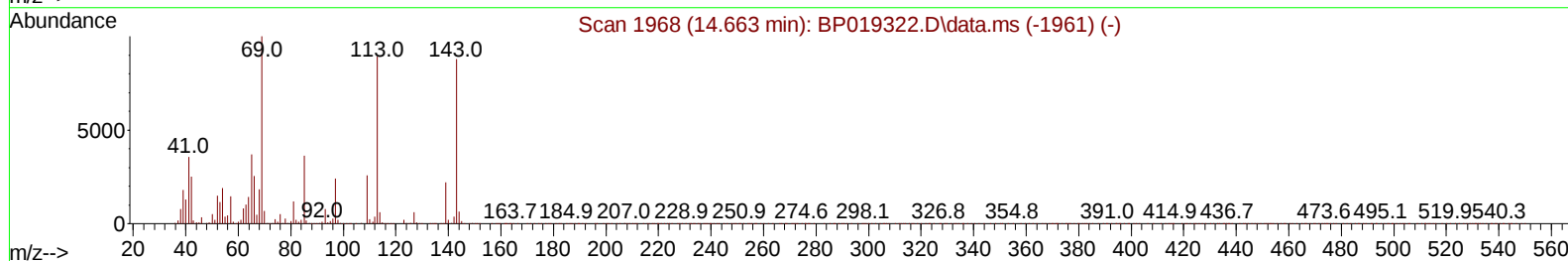
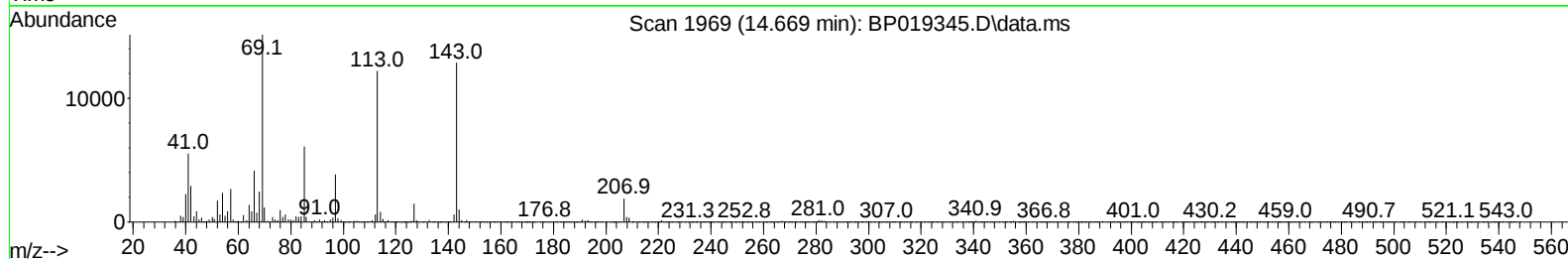
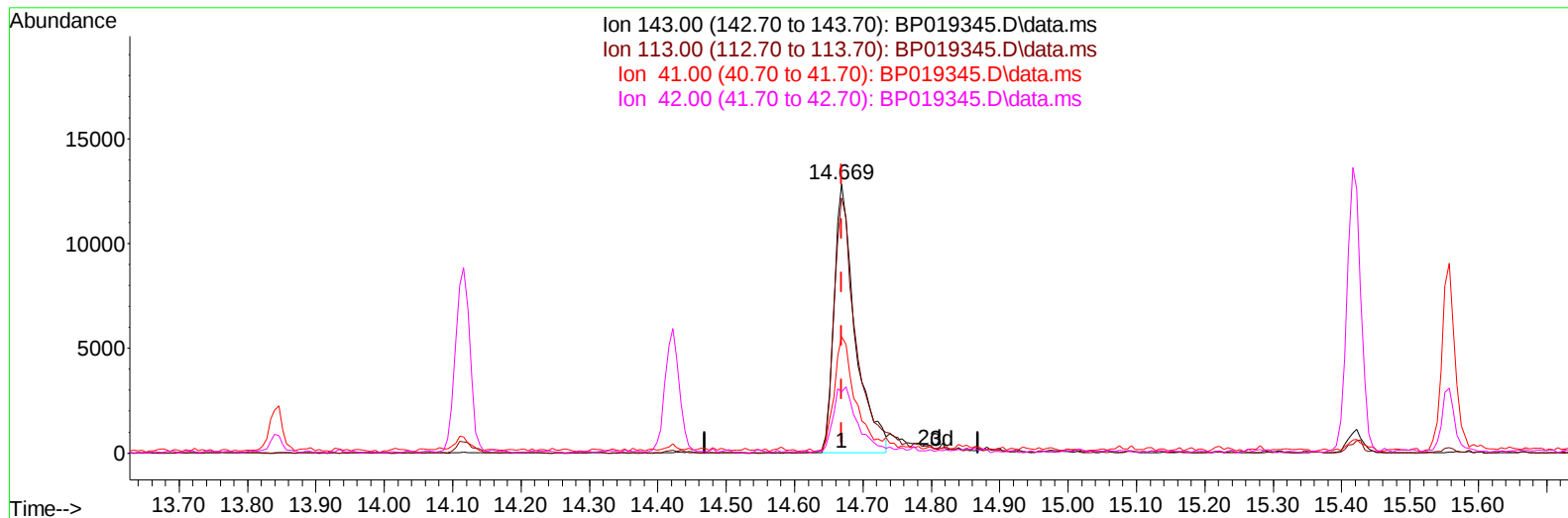
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP011224\
Data File : BP019345.D
Acq On : 12 Jan 2024 23:47
Operator : MA/JU
Sample : P1091-15
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0AZ6

Manual IntegrationsAPPROVED

Quant Time: Jan 13 00:14:57 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: BP019345.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.669min (+ 0.000) 7.53 ng/u1

response 27444

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	98.20	94.86
41.00	39.60	43.26
42.00	25.70	22.68

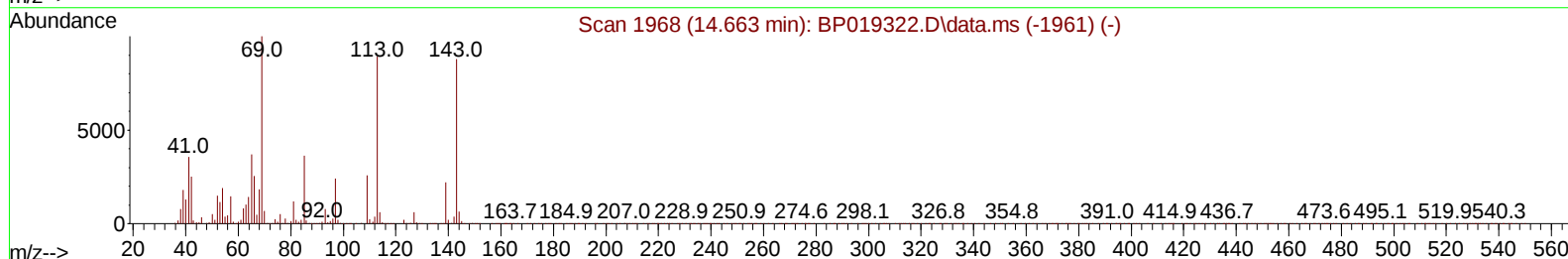
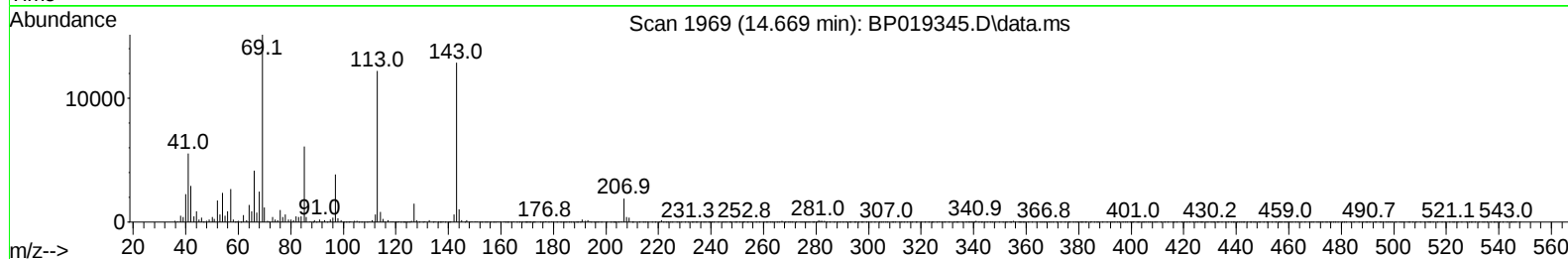
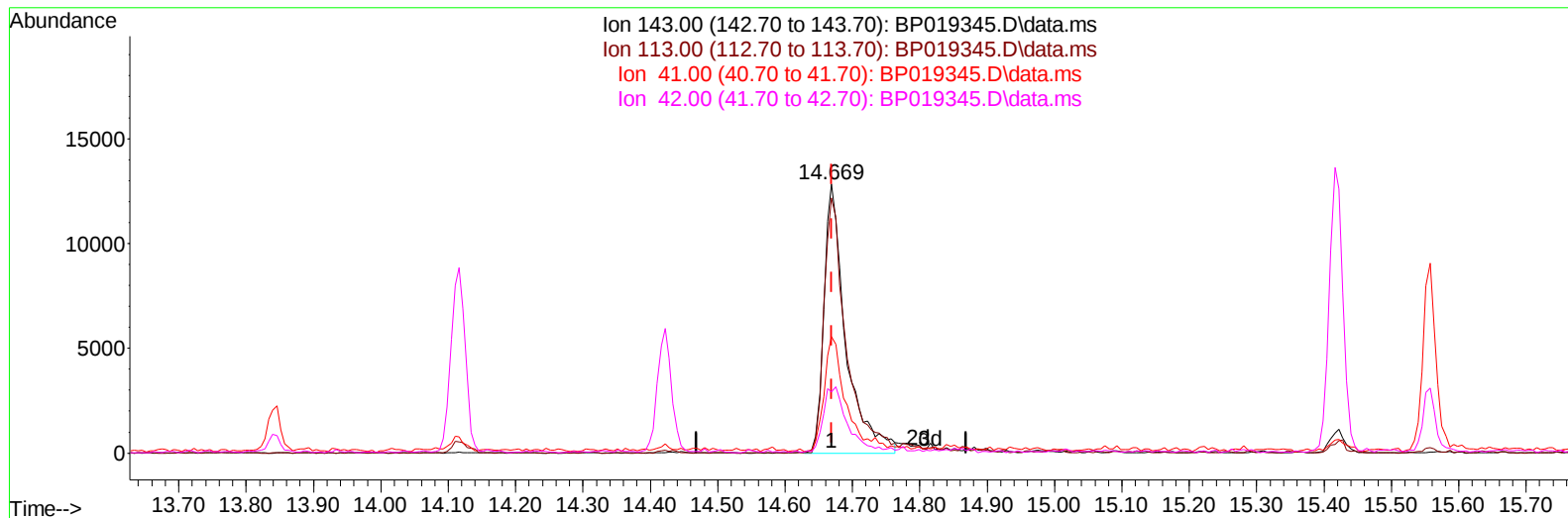
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(54) 4-Nitrophenol-d4 (S)

14.669min (+ 0.000) 7.91 ng/ul m

response 28817

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	98.20	94.86
41.00	39.60	43.26
42.00	25.70	22.68

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.781	152	98902	20.000	ng/ul	0.00
20) Naphthalene-d8	10.575	136	393733	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.422	164	268704	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.181	188	674951	20.000	ng/ul	0.00
79) Chrysene-d12	21.286	240	674812	20.000	ng/ul	0.00
88) Perylene-d12	23.639	264	687973	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.217	96	13823	5.285	ng/uL	0.00
4) Pyridine-d5	3.640	84	92600	12.244	ng/ul	0.00
7) Phenol-d5	6.969	99	80484	8.514	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.116	67	228744	40.456	ng/ul	-0.01
11) 2-Chlorophenol-d4	7.311	132	205068	31.581	ng/ul	0.00
15) 4-Methylphenol-d8	8.505	113	146488	19.533	ng/ul	-0.01
21) Nitrobenzene-d5	8.946	128	134380	40.568	ng/ul	0.00
24) 2-Nitrophenol-d4	9.663	143	155290	40.348	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.204	165	274219	35.266	ng/ul	0.00
31) 4-Chloroaniline-d4	10.722	131	327482	33.341	ng/ul	0.00
46) Dimethylphthalate-d6	13.845	166	987788	41.610	ng/ul	0.00
49) Acenaphthylene-d8	14.116	160	1036314	40.724	ng/ul	0.00
54) 4-Nitrophenol-d4	14.669	143	28817m	7.905	ng/ul	0.00
60) Fluorene-d10	15.422	176	861158	44.308	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.557	200	189058	37.940	ng/ul	0.00
73) Anthracene-d10	17.281	188	1439926	42.775	ng/ul	0.00
81) Pyrene-d10	19.528	212	1848786	42.016	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.486	264	1641354	43.877	ng/ul	0.00

Target Compounds Qvalue

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

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