

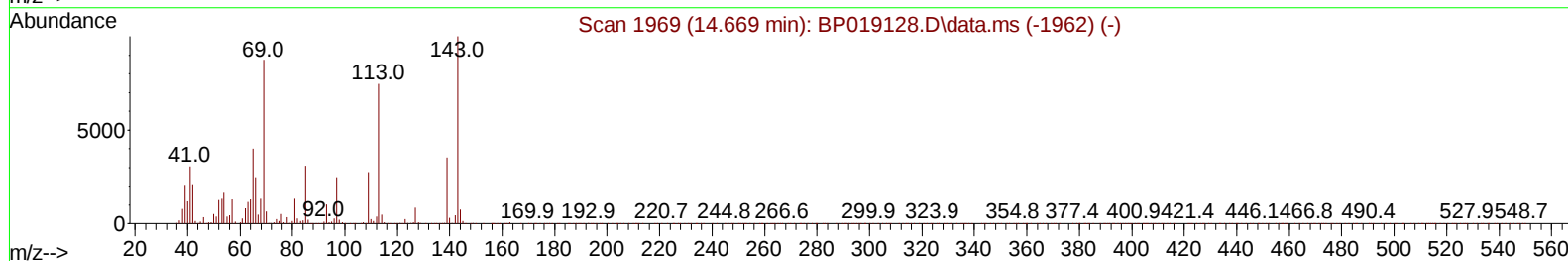
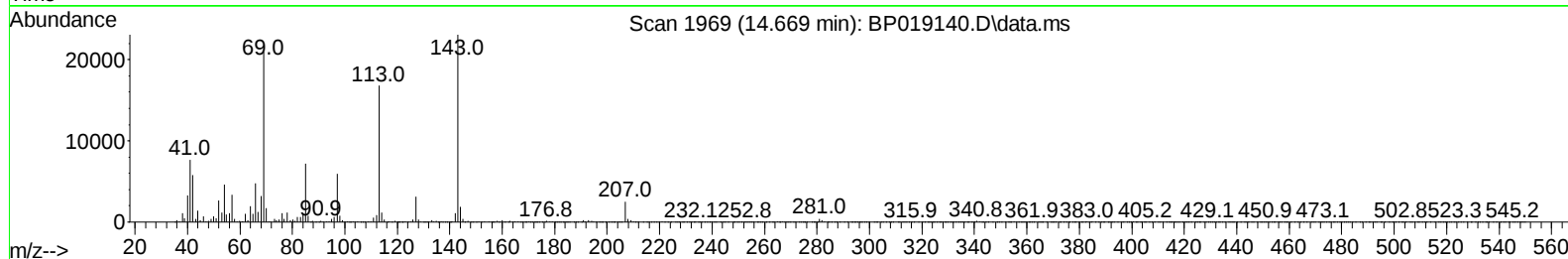
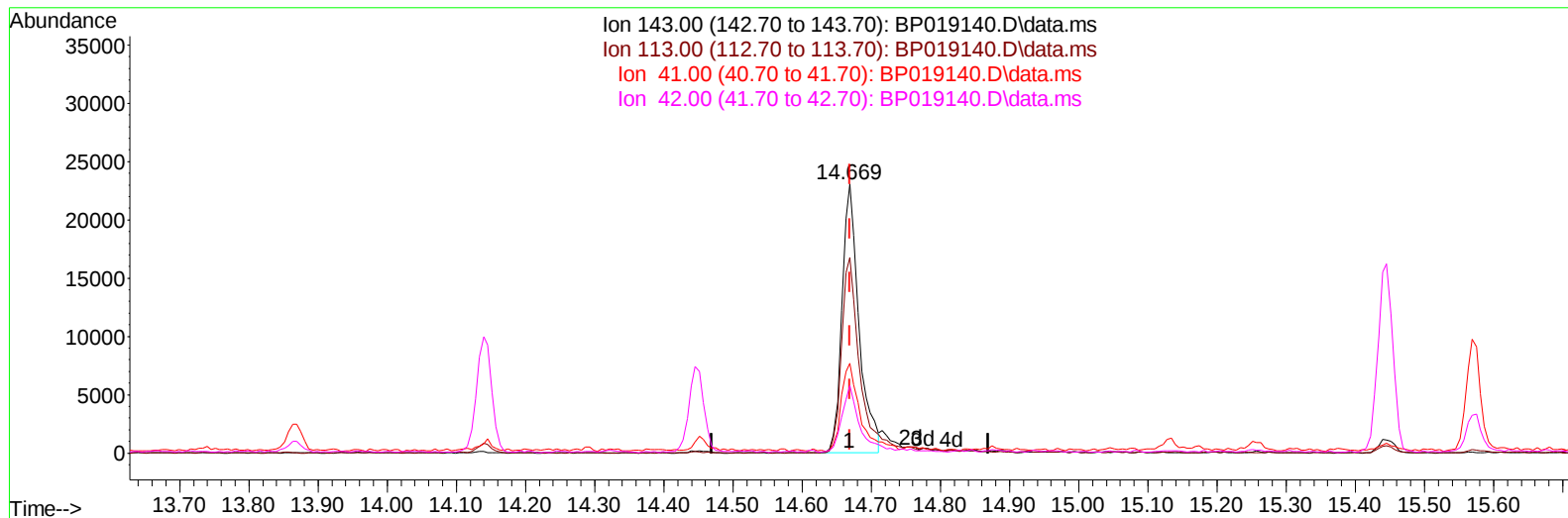
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122823\
Data File : BP019140.D
Acq On : 28 Dec 2023 23:55
Operator : MA/JU
Sample : 06021-18
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BH800

Manual IntegrationsAPPROVED

Quant Time: Dec 29 00:22:25 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP122723.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Thu Dec 28 03:15:29 2023
Response via : Initial Calibration

Reviewed By :Yogesh Patel 12/29/2023
Supervised By :mohammad ahmed 01/01/2024



TIC: BP019140.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.669min (-0.000) 6.37 ng/u1

response 39305

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	73.50	72.74
41.00	33.40	33.29
42.00	23.00	25.14

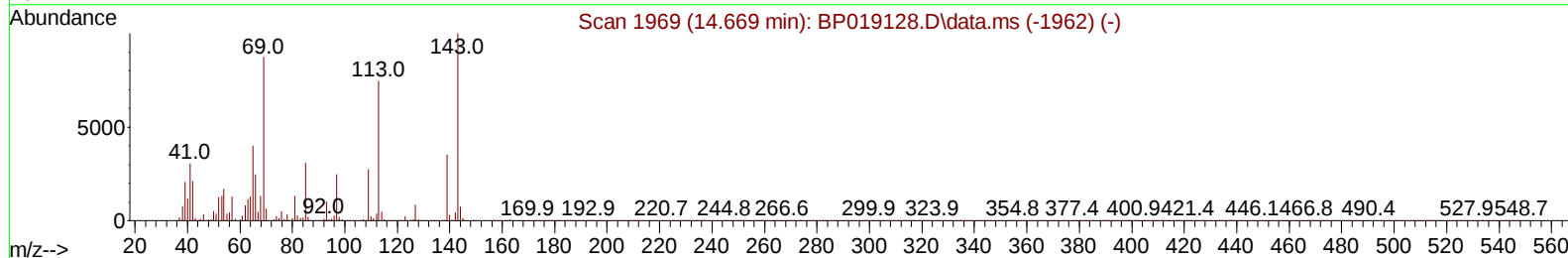
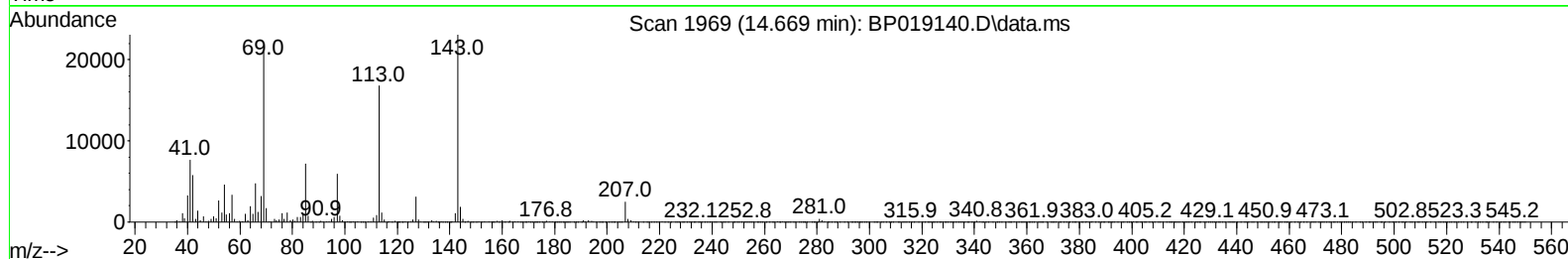
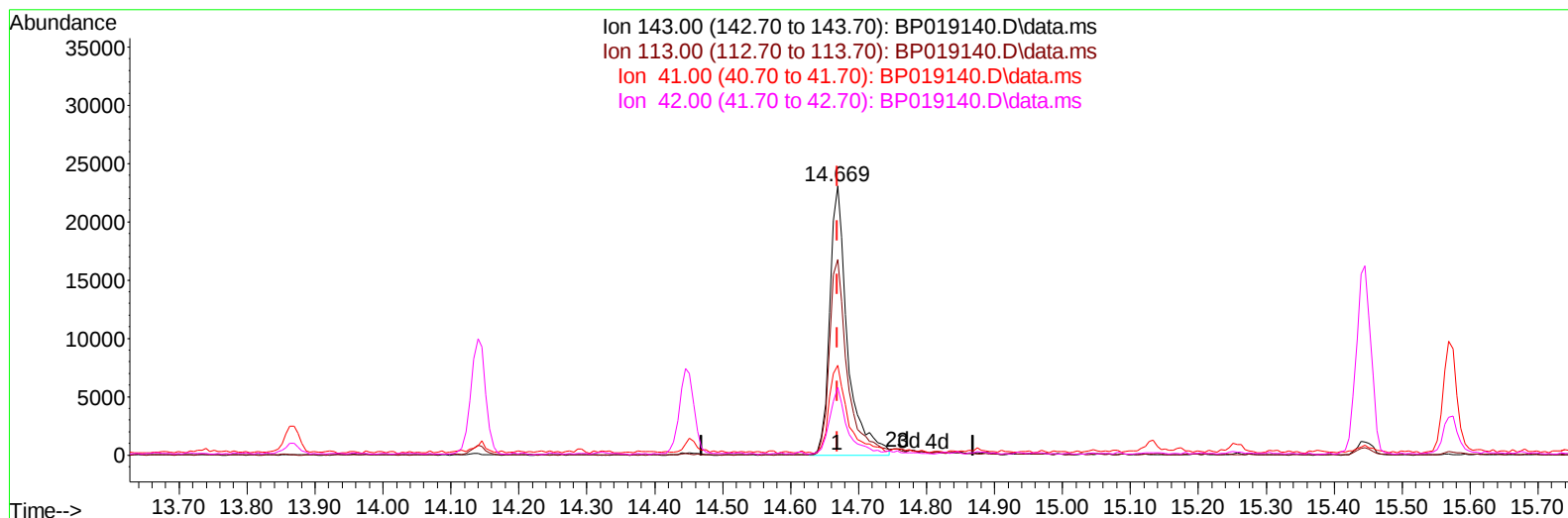
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TIC: BP019140.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.669min (-0.000) 6.79 ng/u1 m

response 41897

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	73.50	72.74
41.00	33.40	33.29
42.00	23.00	25.14

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Quant Time: Dec 29 00:23:12 2023
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 QLast Update : Thu Dec 28 03:15:29 2023
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.805	152	176598	20.000	ng/ul	0.00
20) Naphthalene-d8	10.598	136	667734	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.451	164	413417	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.204	188	928698	20.000	ng/ul	0.00
79) Chrysene-d12	21.310	240	1013663	20.000	ng/ul	0.00
88) Perylene-d12	23.674	264	1210630	20.000	ng/ul	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.234	96	23863	4.615	ng/uL	0.00
4) Pyridine-d5	3.658	84	78970	6.085	ng/ul	0.00
7) Phenol-d5	6.981	99	137620	8.690	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.146	67	364718	39.808	ng/ul	0.00
11) 2-Chlorophenol-d4	7.334	132	375369	29.401	ng/ul	0.00
15) 4-Methylphenol-d8	8.522	113	235905	19.133	ng/ul	0.00
21) Nitrobenzene-d5	8.969	128	226036	39.160	ng/ul	0.00
24) 2-Nitrophenol-d4	9.693	143	243824	39.085	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.228	165	436500	34.651	ng/ul	0.00
31) 4-Chloroaniline-d4	10.746	131	337225	21.764	ng/ul	0.00
46) Dimethylphthalate-d6	13.869	166	1399461	38.324	ng/ul	0.00
49) Acenaphthylene-d8	14.139	160	1507361	38.043	ng/ul	0.00
54) 4-Nitrophenol-d4	14.669	143	41897m	6.795	ng/ul	0.00
60) Fluorene-d10	15.445	176	1212274	38.428	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.569	200	210309	34.543	ng/ul	-0.01
73) Anthracene-d10	17.304	188	1973174	40.684	ng/ul	0.00
81) Pyrene-d10	19.551	212	2484125	40.211	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.521	264	2843609	40.864	ng/ul	-0.01
Target Compounds						
					Qvalue	
35) 4-Chloro-3-methylphenol	11.904	107	17622	1.458	ng/ul	98

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

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