

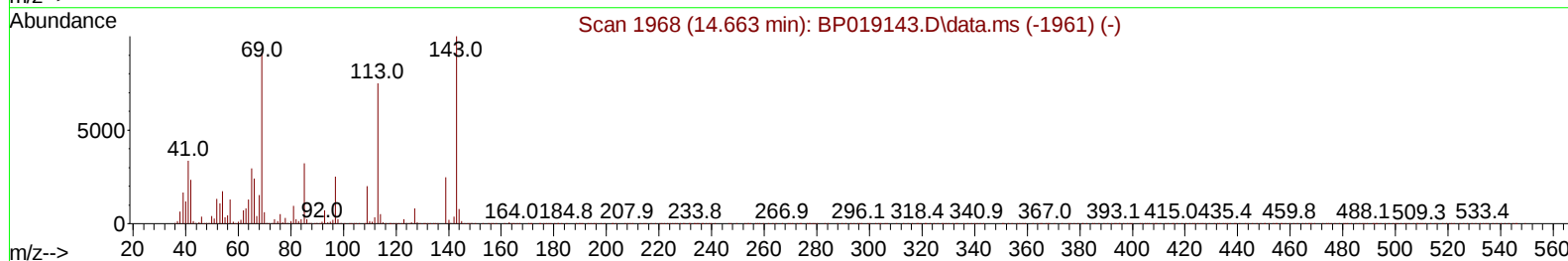
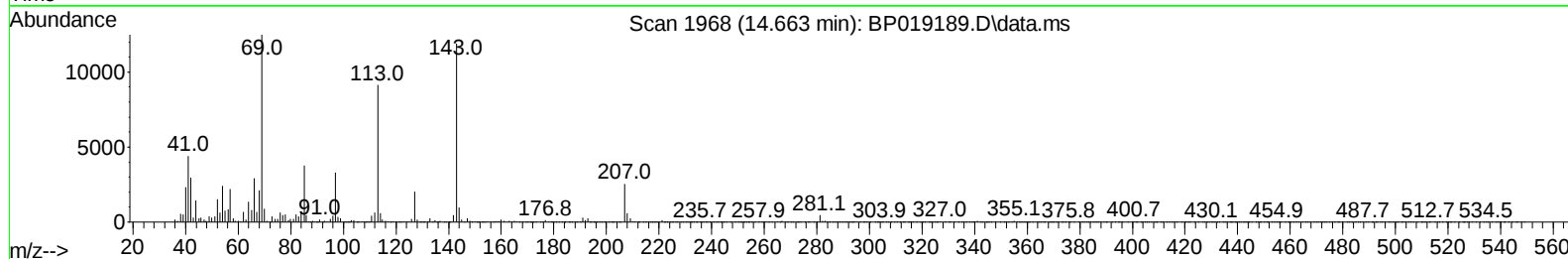
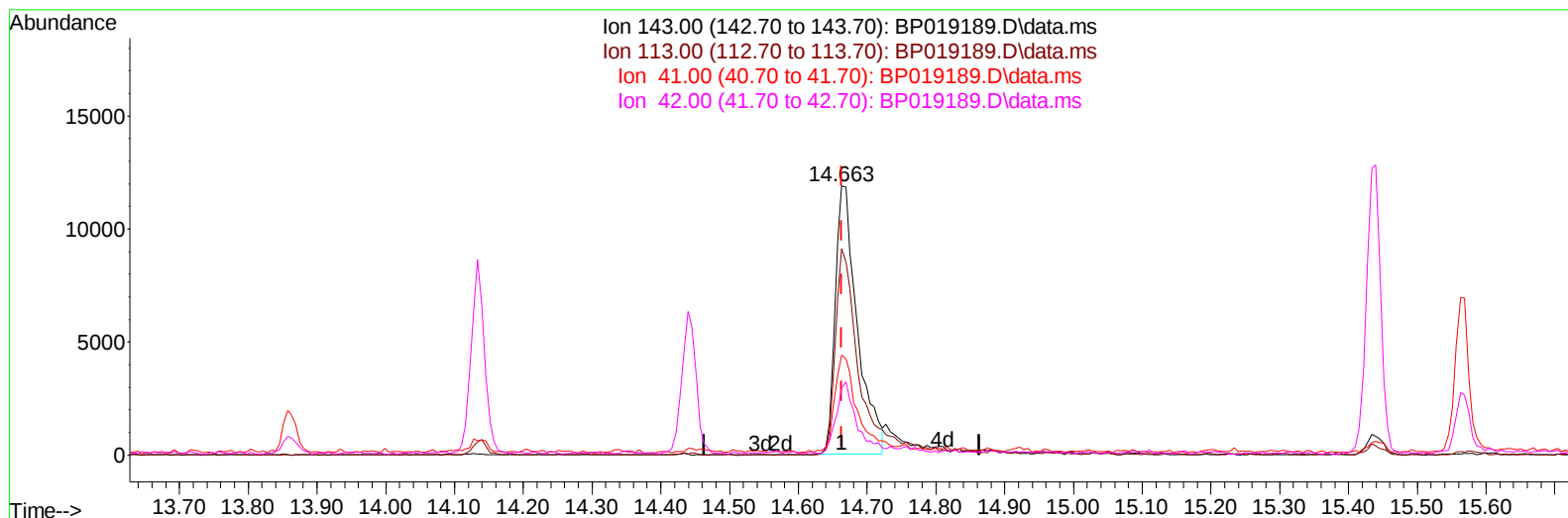
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122823\
Data File : BP019189.D
Acq On : 30 Dec 2023 04:55
Operator : MA/JU
Sample : 06020-11
Misc :
ALS Vial : 68 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BH7X6

Manual IntegrationsAPPROVED

Quant Time: Jan 01 00:19:39 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP122723.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri Dec 29 03:48:11 2023
Response via : Initial Calibration

Reviewed By :Yogesh Patel 01/01/2024
Supervised By :mohammad ahmed 01/01/2024



TIC: BP019189.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.663min (-0.000) 5.23 ng/u1

response 26356

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	73.50	76.85
41.00	33.40	37.11
42.00	23.00	24.87

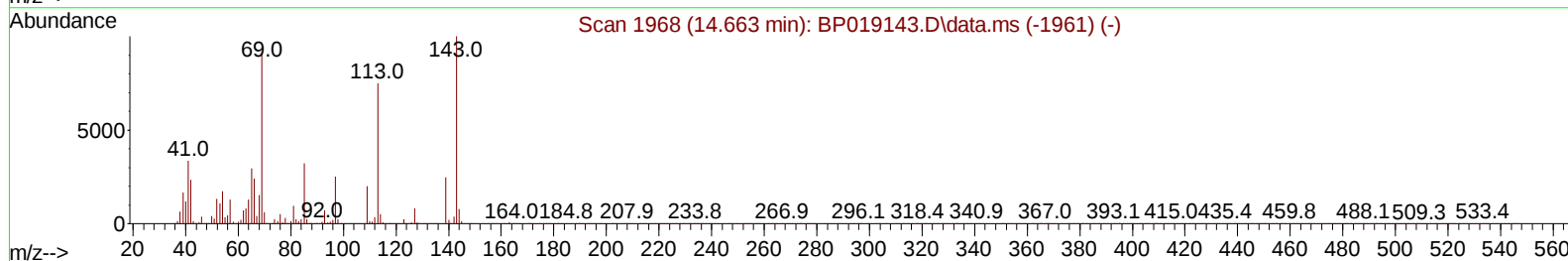
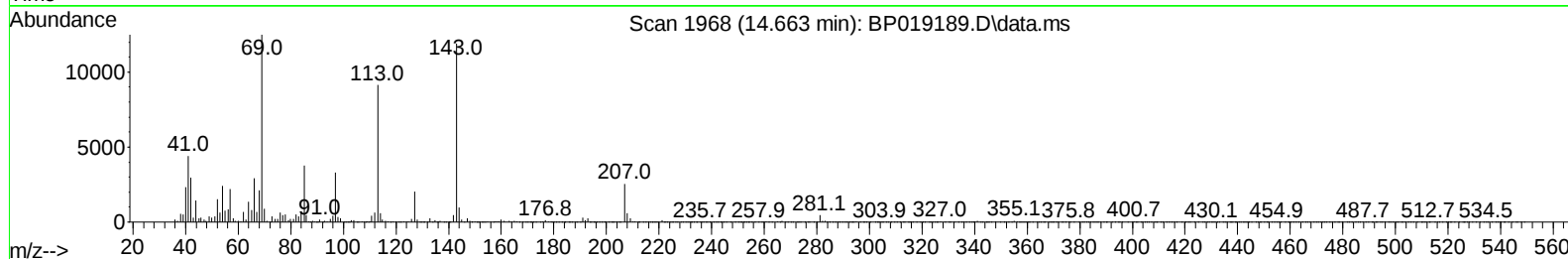
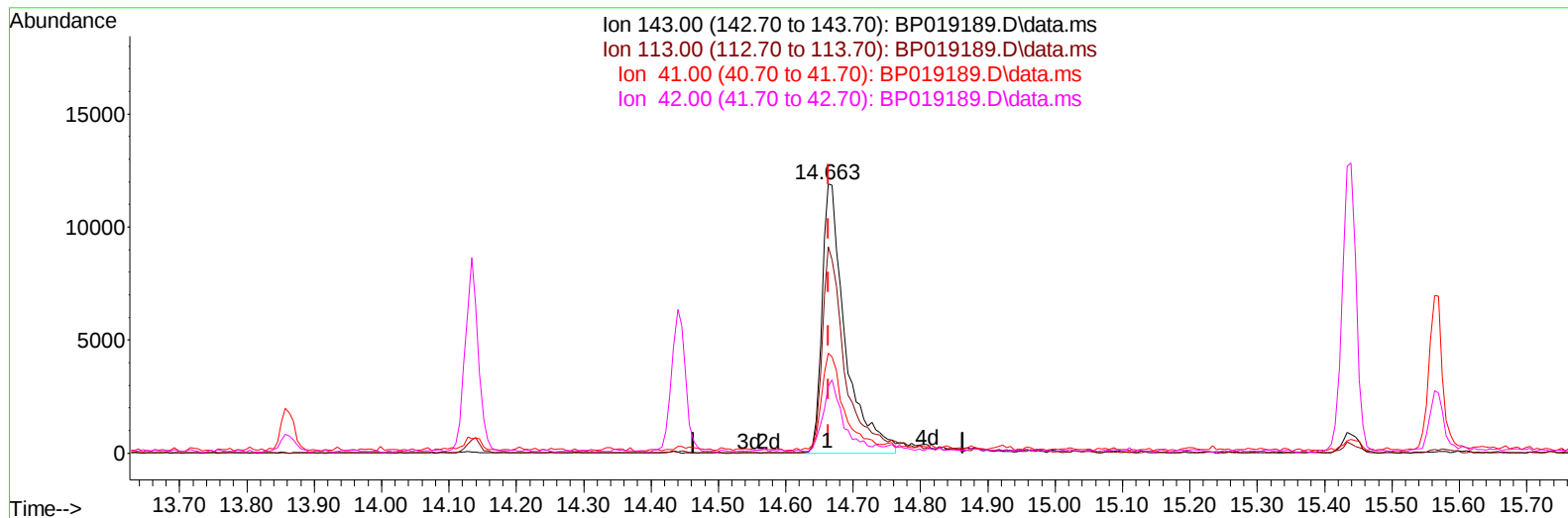
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TIC: BP019189.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.663min (-0.000) 5.67 ng/ul m

response 28538

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	73.50	76.85
41.00	33.40	37.11
42.00	23.00	24.87

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Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 01/01/2024
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Quant Time: Jan 01 00:24:43 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP122723.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Dec 29 03:48:11 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.799	152	160074	20.000	ng/ul	0.00
20) Naphthalene-d8	10.593	136	589851	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.440	164	337612	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.198	188	760379	20.000	ng/ul	0.00
79) Chrysene-d12	21.304	240	844205	20.000	ng/ul	0.00
88) Perylene-d12	23.663	264	1012774	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.234	96	20684	4.413	ng/uL	0.00
4) Pyridine-d5	3.658	84	117339	9.974	ng/ul	0.00
7) Phenol-d5	6.975	99	107434	7.484	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.140	67	280857	33.819	ng/ul	0.00
11) 2-Chlorophenol-d4	7.328	132	288103	24.895	ng/ul	0.00
15) 4-Methylphenol-d8	8.516	113	179803	16.088	ng/ul	0.00
21) Nitrobenzene-d5	8.963	128	161130	31.601	ng/ul	0.00
24) 2-Nitrophenol-d4	9.681	143	162750	29.534	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.222	165	303776	27.299	ng/ul	0.00
31) 4-Chloroaniline-d4	10.740	131	377705	27.595	ng/ul	0.00
46) Dimethylphthalate-d6	13.863	166	1010943	33.901	ng/ul	0.00
49) Acenaphthylene-d8	14.134	160	1064105	32.886	ng/ul	0.00
54) 4-Nitrophenol-d4	14.663	143	28538m	5.667	ng/ul	0.00
60) Fluorene-d10	15.439	176	899089	34.899	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.569	200	142496	28.586	ng/ul	0.00
73) Anthracene-d10	17.298	188	1581752	39.833	ng/ul	0.00
81) Pyrene-d10	19.545	212	2005467	38.979	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.515	264	2334441	40.101	ng/ul	0.00

Target Compounds Qvalue

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

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