

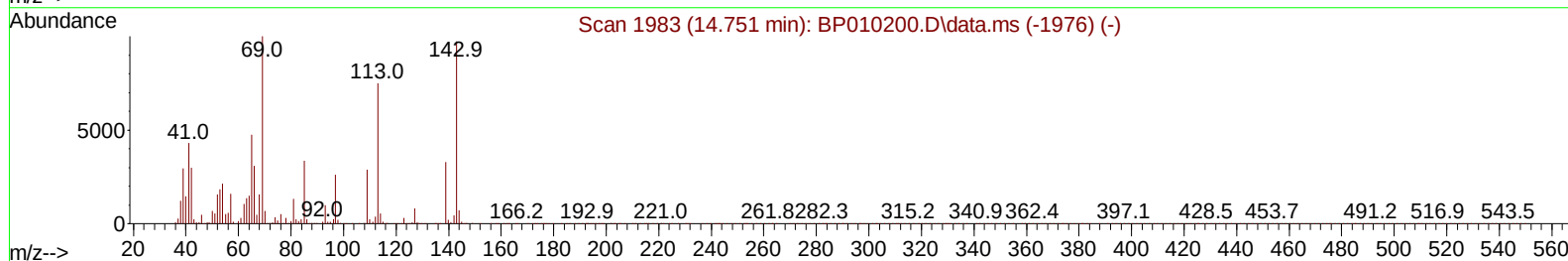
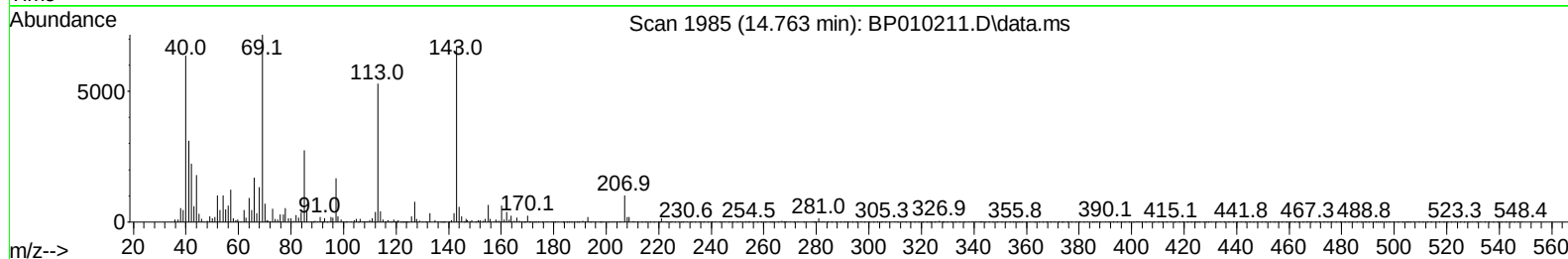
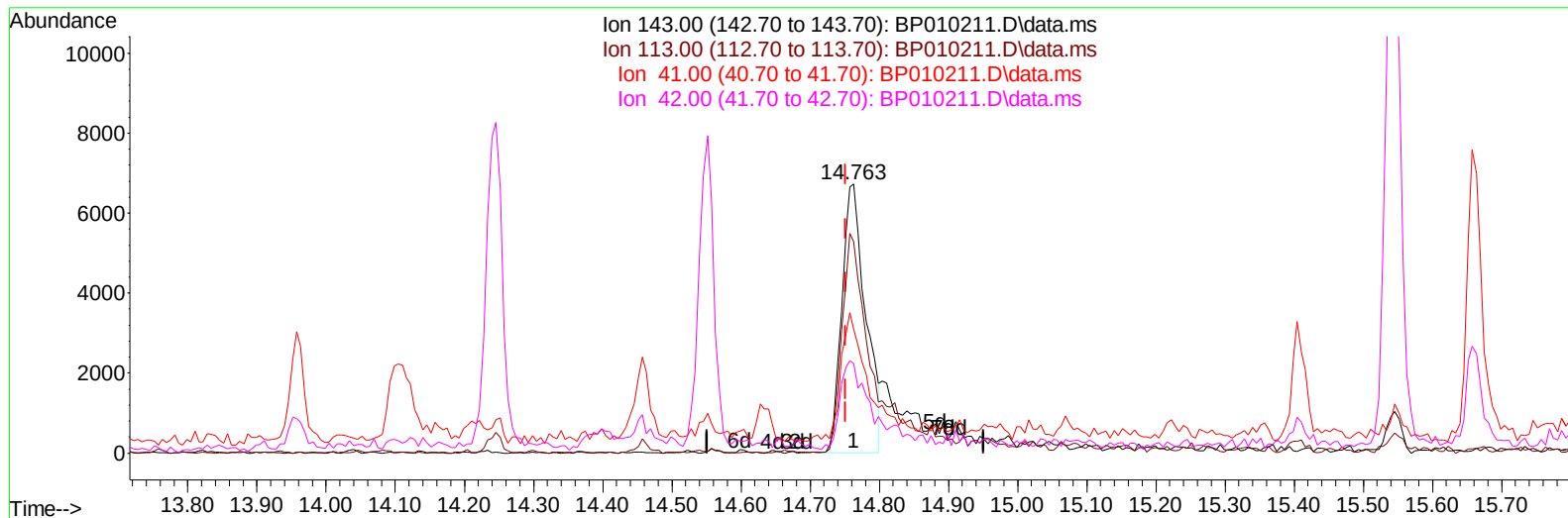
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP050222\
Data File : BP010211.D
Acq On : 03 May 2022 17:16
Operator : CG/JU
Sample : N2615-19
Misc :
ALS Vial : 47 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BH0D9

Manual IntegrationsAPPROVED

Quant Time: May 04 01:10:43 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP042122.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue May 03 00:58:32 2022
Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 05/04/2022
Supervised By :Yogesh Patel 05/05/2022



TIC: BP010211.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.763min (+ 0.012) 3.18 ng/ul

response 15979

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	132.70	78.58#
41.00	23.20	46.22#
42.00	18.00	33.39#

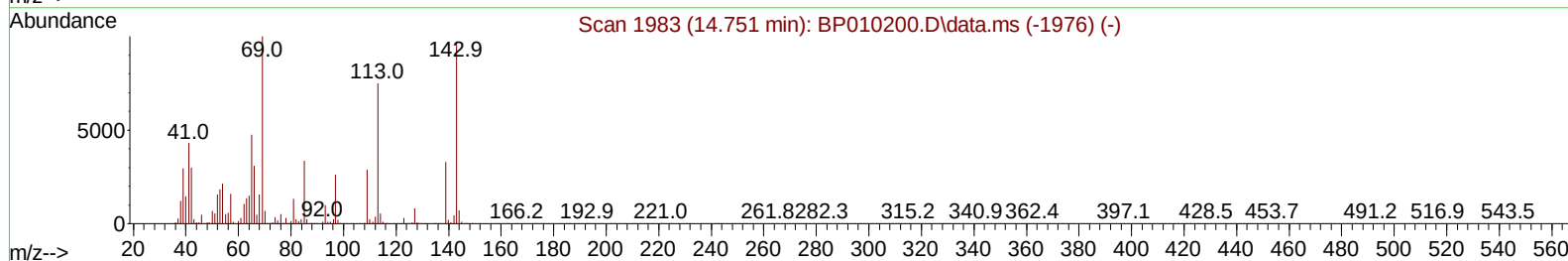
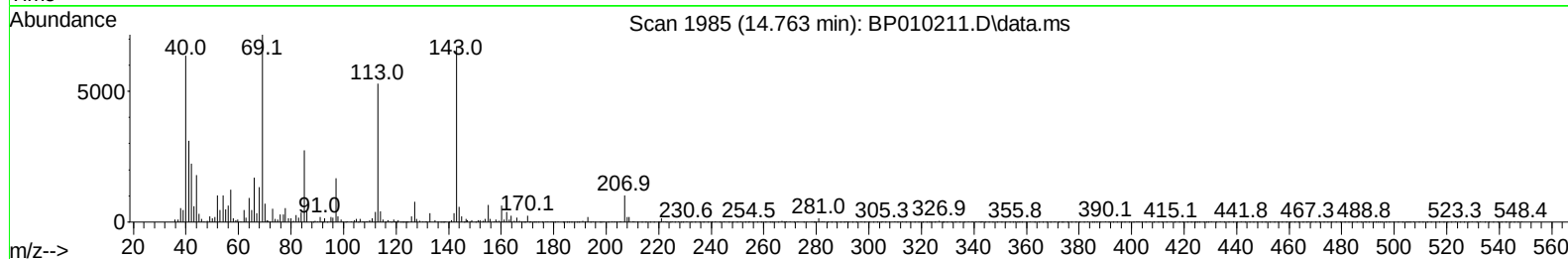
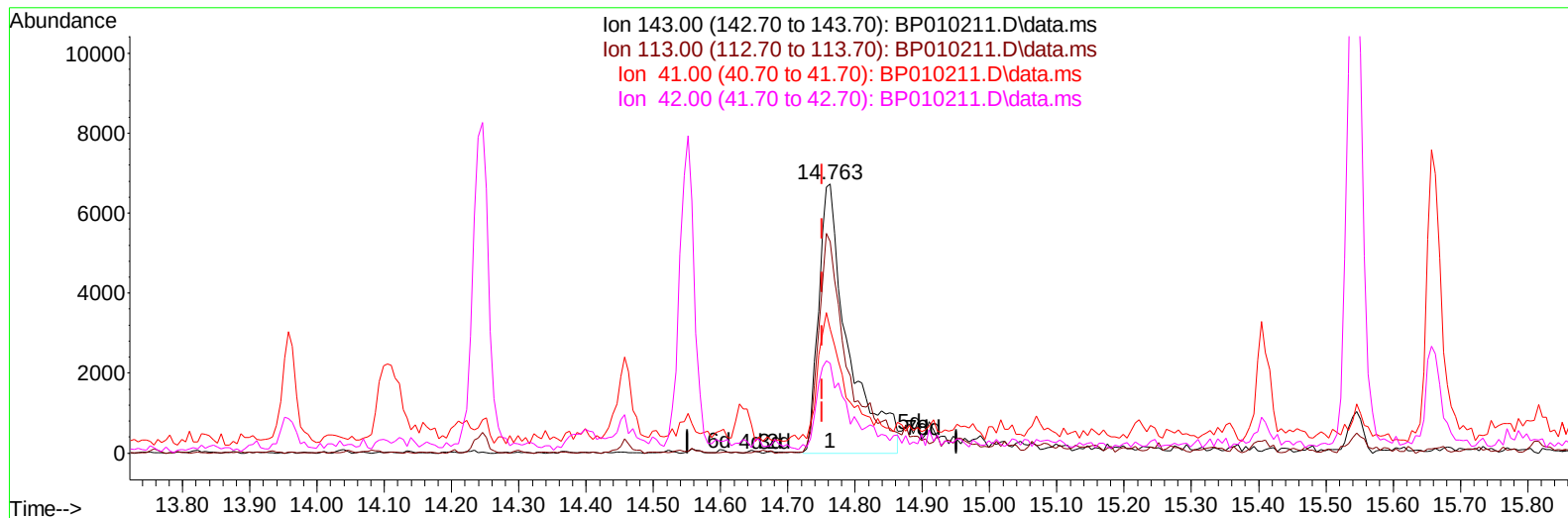
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TIC: BP010211.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.763min (+ 0.012) 4.06 ng/ul m

response 20398

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	132.70	78.58#
41.00	23.20	46.22#
42.00	18.00	33.39#

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.922	152	135367	20.000	ng/ul	0.00
20) Naphthalene-d8	10.722	136	553624	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.551	164	350608	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.304	188	717407	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.386	240	655255	20.000	ng/ul	# 0.00
88) Perylene-d12	23.827	264	659402	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.358	96	15498	4.992	ng/uL	0.00
4) Pyridine-d5	3.781	84	81524	9.335	ng/ul	0.00
7) Phenol-d5	7.081	99	70882	5.757	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.246	67	245791	33.268	ng/ul	0.00
11) 2-Chlorophenol-d4	7.452	132	217915	23.897	ng/ul	0.00
15) 4-Methylphenol-d8	8.628	113	136159	13.347	ng/ul	0.00
21) Nitrobenzene-d5	9.081	128	144798	32.885	ng/ul	0.00
24) 2-Nitrophenol-d4	9.805	143	152636	31.740	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.346	165	239174	25.548	ng/ul	0.00
31) 4-Chloroaniline-d4	10.858	131	324239	24.155	ng/ul	0.00
46) Dimethylphthalate-d6	13.957	166	913116	33.480	ng/ul	0.00
49) Acenaphthylene-d8	14.246	160	1045506	31.674	ng/ul	0.00
54) 4-Nitrophenol-d4	14.763	143	20398m	4.063	ng/ul	0.01
60) Fluorene-d10	15.545	176	799171	34.064	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.663	200	114294	25.196	ng/ul	0.00
73) Anthracene-d10	17.404	188	1342167	38.671	ng/ul	0.00
81) Pyrene-d10	19.634	212	1550263	41.602	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.669	264	1432316	41.433	ng/ul	-0.01

Target Compounds Qvalue

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

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