

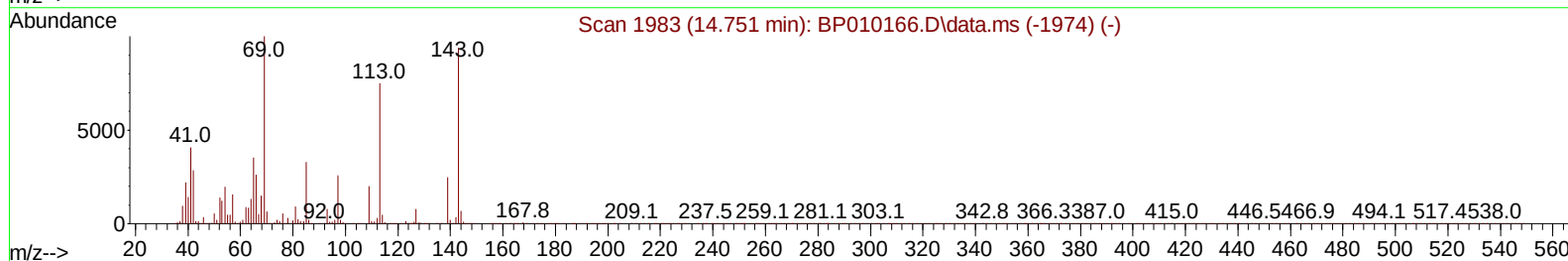
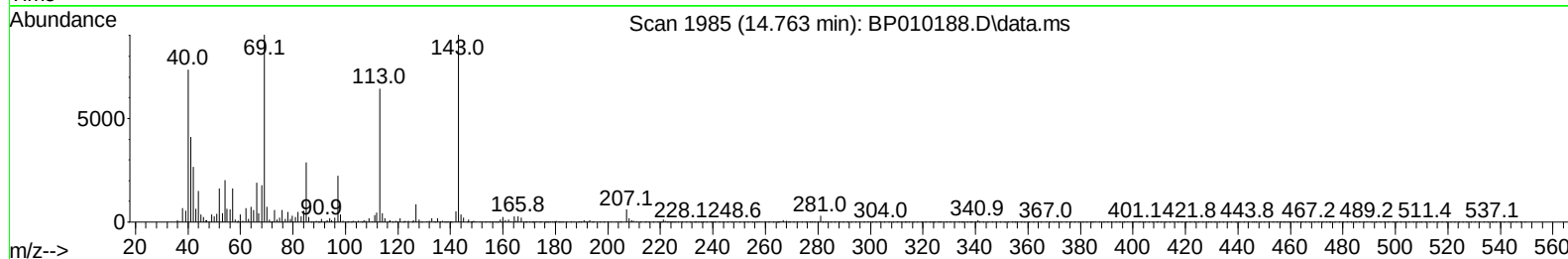
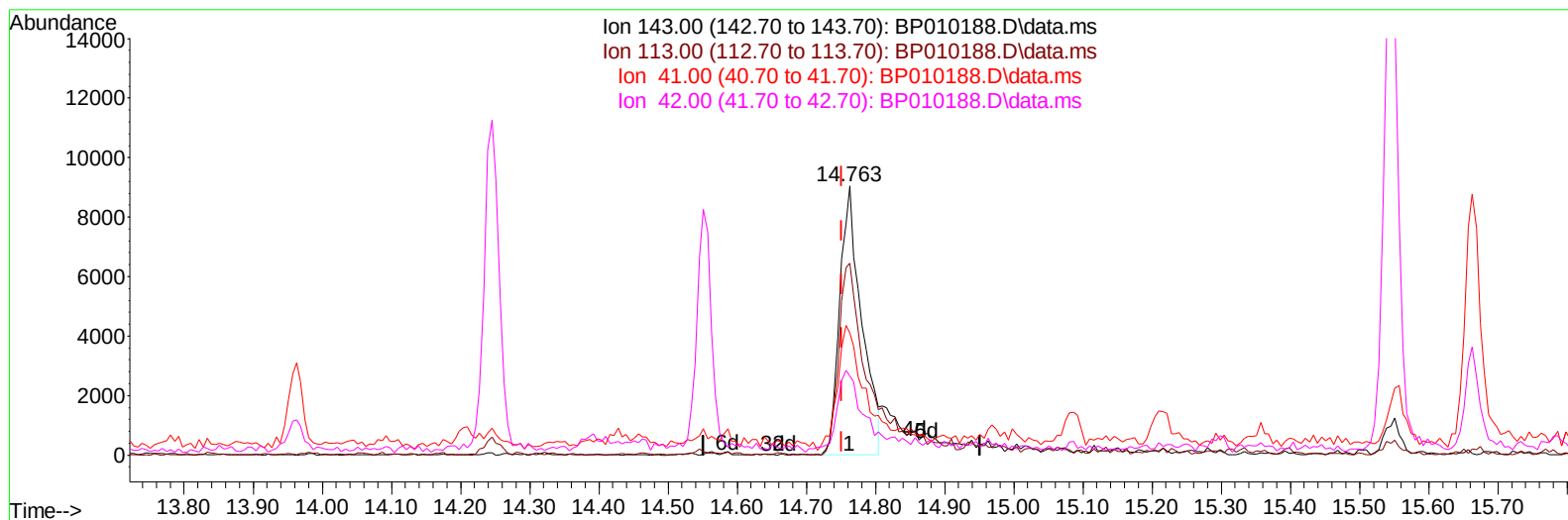
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP050222\
Data File : BP010188.D
Acq On : 03 May 2022 00:16
Operator : CG/JU
Sample : N2615-16
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BH0C8

Manual IntegrationsAPPROVED

Quant Time: May 03 02:53:19 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 05/03/2022
Supervised By :Yogesh Patel 05/05/2022



TIC: BP010188.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.763min (+ 0.011) 4.08 ng/u1

response 19894

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	132.70	71.23#
41.00	23.20	45.38#
42.00	18.00	29.44#

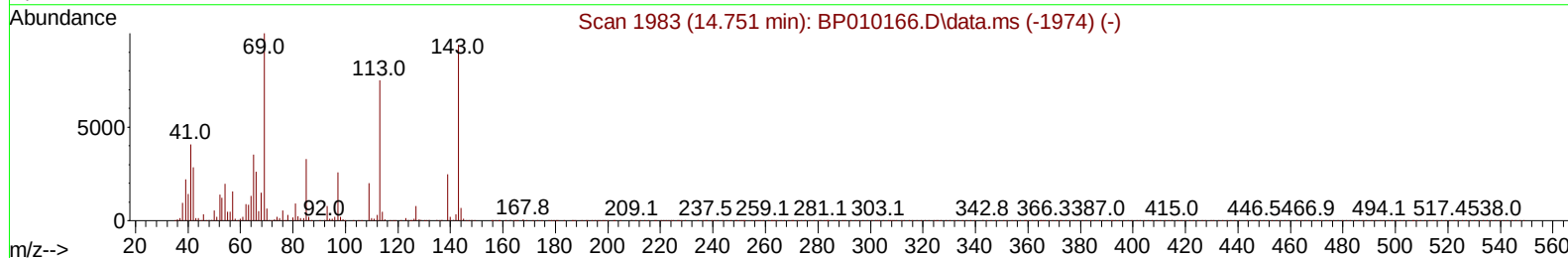
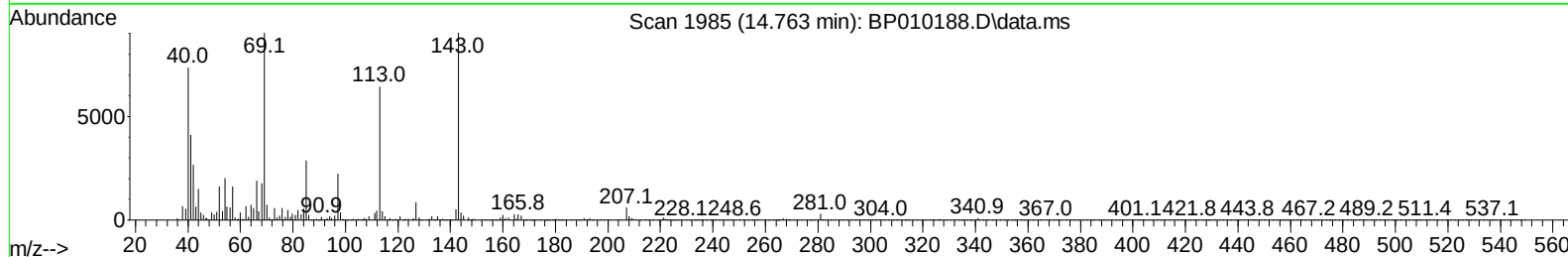
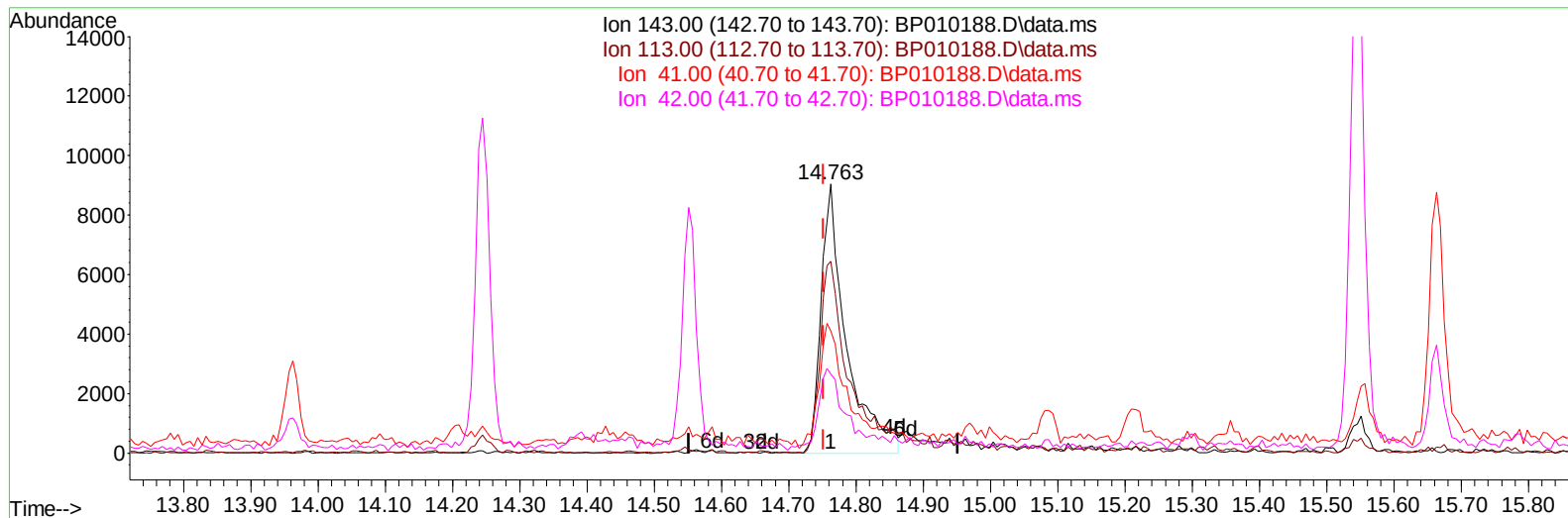
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TIC: BP010188.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.763min (+ 0.011) 4.84 ng/ul m

response 23584

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	132.70	71.23#
41.00	23.20	45.38#
42.00	18.00	29.44#

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Instrument :
 BNA_P
 ClientSampleId :
 BH0C8

Manual IntegrationsAPPROVED

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.922	152	135239	20.000	ng/ul	0.00
20) Naphthalene-d8	10.728	136	565892	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.551	164	340479	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.304	188	623096	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.386	240	524427	20.000	ng/ul	# 0.00
88) Perylene-d12	23.833	264	520886	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.357	96	17290	5.575	ng/uL	0.00
4) Pyridine-d5	3.781	84	100498	11.519	ng/ul	0.00
7) Phenol-d5	7.087	99	85145	6.922	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.245	67	301804	40.888	ng/ul	0.00
11) 2-Chlorophenol-d4	7.451	132	270181	29.657	ng/ul	0.00
15) 4-Methylphenol-d8	8.628	113	173400	17.013	ng/ul	0.00
21) Nitrobenzene-d5	9.081	128	186562	41.451	ng/ul	0.00
24) 2-Nitrophenol-d4	9.804	143	194939	39.658	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.345	165	307823	32.168	ng/ul	0.00
31) 4-Chloroaniline-d4	10.863	131	415727	30.299	ng/ul	0.00
46) Dimethylphthalate-d6	13.963	166	1124863	42.471	ng/ul	0.00
49) Acenaphthylene-d8	14.245	160	1272279	39.691	ng/ul	0.00
54) 4-Nitrophenol-d4	14.763	143	23584m	4.837	ng/ul	0.01
60) Fluorene-d10	15.545	176	935420	41.058	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.663	200	139137	35.315	ng/ul	0.00
73) Anthracene-d10	17.404	188	1351030	44.818	ng/ul	0.00
81) Pyrene-d10	19.633	212	1512171	50.703	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.674	264	1329273	48.677	ng/ul	0.00

Target Compounds Qvalue

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

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