

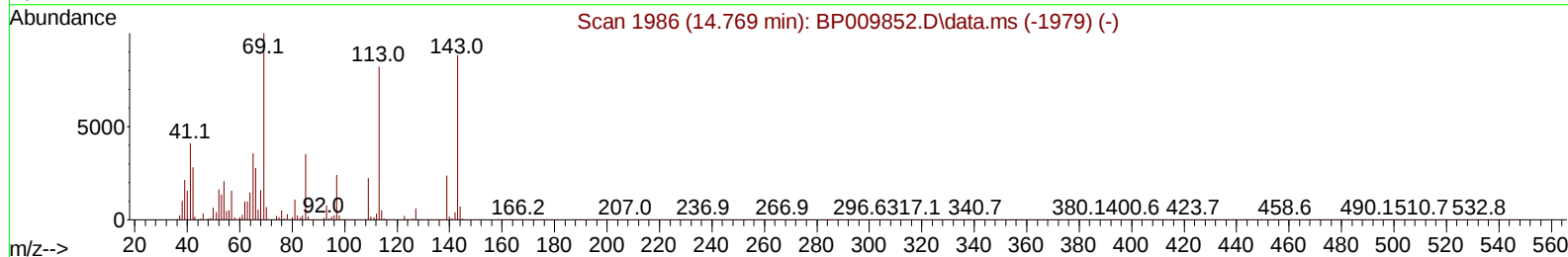
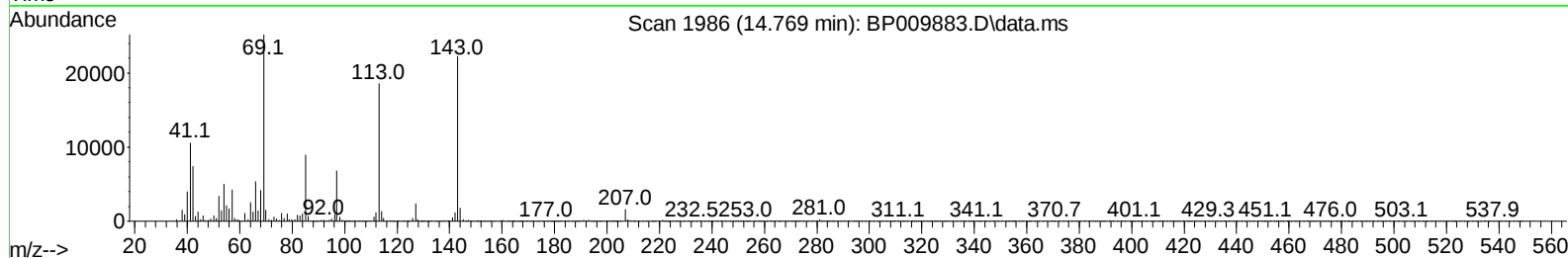
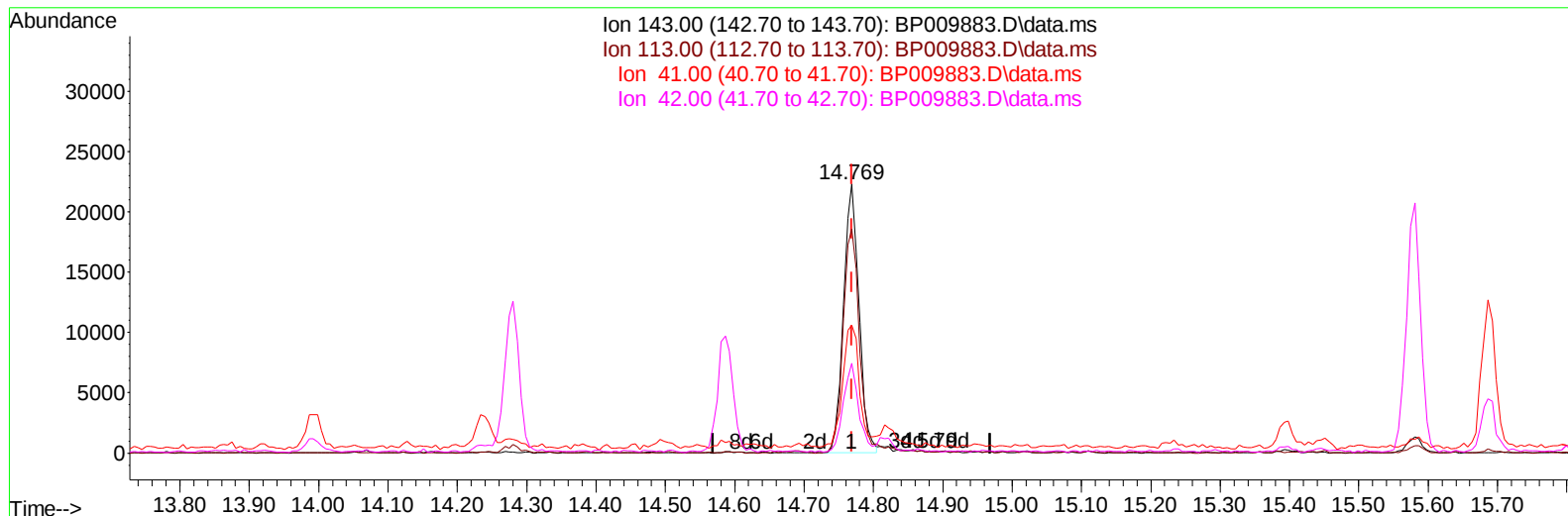
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041522\
Data File : BP009883.D
Acq On : 16 Apr 2022 03:58
Operator : CG/JU
Sample : N2421-04
Misc :
ALS Vial : 34 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0J74

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 04/18/2022
Supervised By :Yogesh Patel 04/18/2022

Quant Time: Apr 18 01:08:19 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP040722.M
Quant Title : SVOA CALIBRATION
QLast Update : Mon Apr 18 00:59:30 2022
Response via : Initial Calibration



TIC: BP009883.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.769min (0.000) 5.69 ng/u1

response 34242

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	132.70	83.24#
41.00	23.20	47.46#
42.00	18.00	33.40#

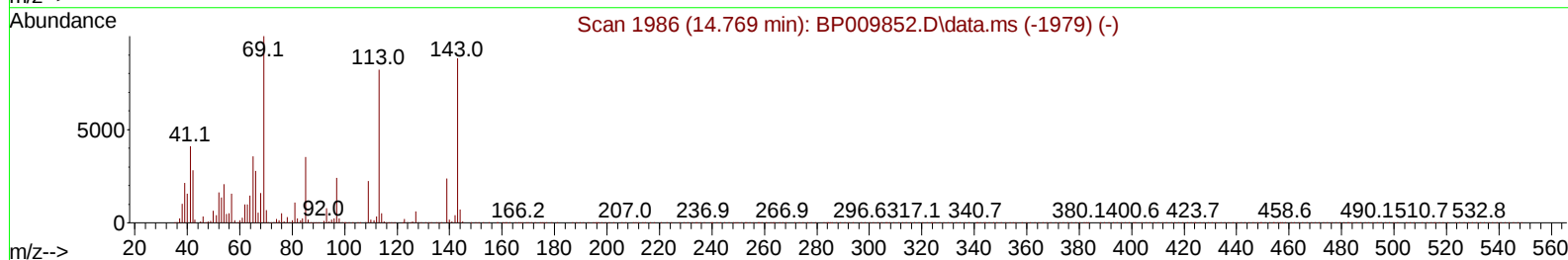
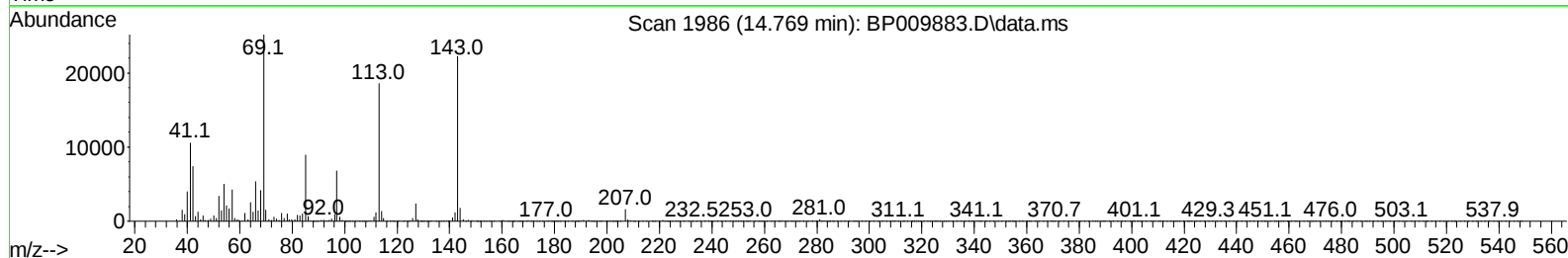
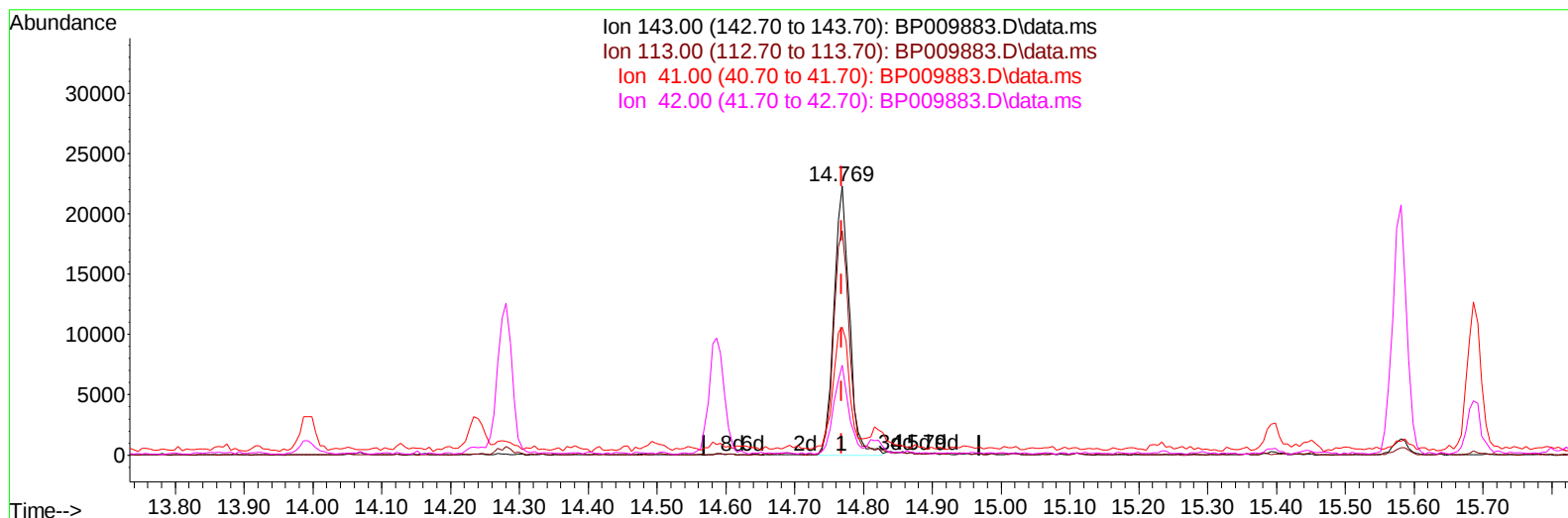
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TIC: BP009883.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.769min (0.000) 5.80 ng/ul m

response 34943

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	132.70	83.24#
41.00	23.20	47.46#
42.00	18.00	33.40#

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.957	152	131900	20.000	ng/ul	0.00
20) Naphthalene-d8	10.763	136	586801	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.587	164	422718	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.339	188	956216	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.416	240	1003064	20.000	ng/ul	# 0.00
88) Perylene-d12	23.880	264	920433	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.381	96	12665	4.285	ng/uL	0.00
4) Pyridine-d5	3.811	84	19720	2.420	ng/ul	0.01
7) Phenol-d5	7.110	99	70611	6.098	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.281	67	242266	35.142	ng/ul	0.00
11) 2-Chlorophenol-d4	7.487	132	219443	25.122	ng/ul	0.00
15) 4-Methylphenol-d8	8.657	113	154670	16.000	ng/ul	0.00
21) Nitrobenzene-d5	9.116	128	161930	38.267	ng/ul	0.00
24) 2-Nitrophenol-d4	9.840	143	167668	36.141	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.375	165	285486	29.407	ng/ul	0.00
31) 4-Chloroaniline-d4	10.893	131	252795	18.283	ng/ul	0.00
46) Dimethylphthalate-d6	13.992	166	1253709	37.757	ng/ul	0.00
49) Acenaphthylene-d8	14.281	160	1350121	34.320	ng/ul	0.00
54) 4-Nitrophenol-d4	14.769	143	34943m	5.803	ng/ul	0.00
60) Fluorene-d10	15.581	176	1076573	38.541	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.686	200	194239	34.290	ng/ul	0.00
73) Anthracene-d10	17.439	188	1853637	40.488	ng/ul	0.00
81) Pyrene-d10	19.669	212	2267879	41.398	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.727	264	2057313	42.921	ng/ul	0.00

Target Compounds Qvalue

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

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