

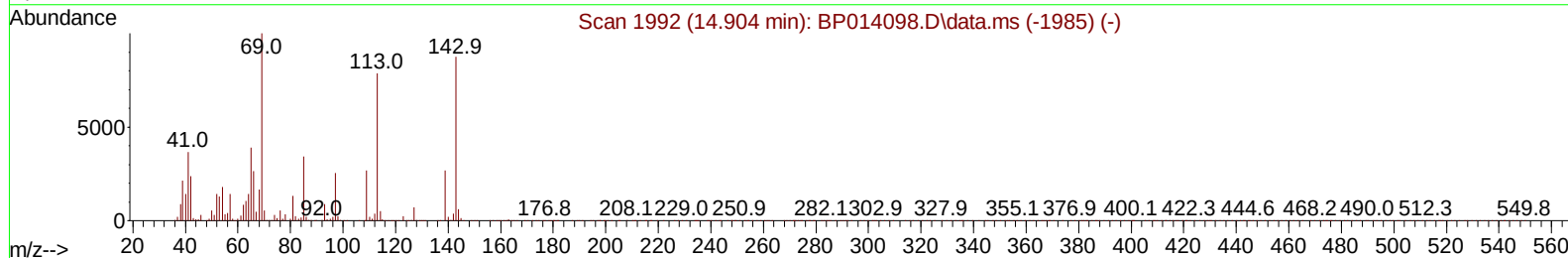
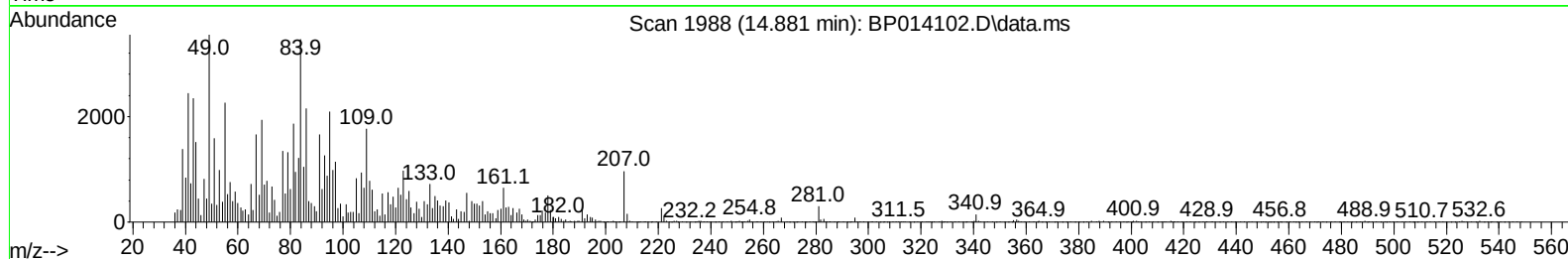
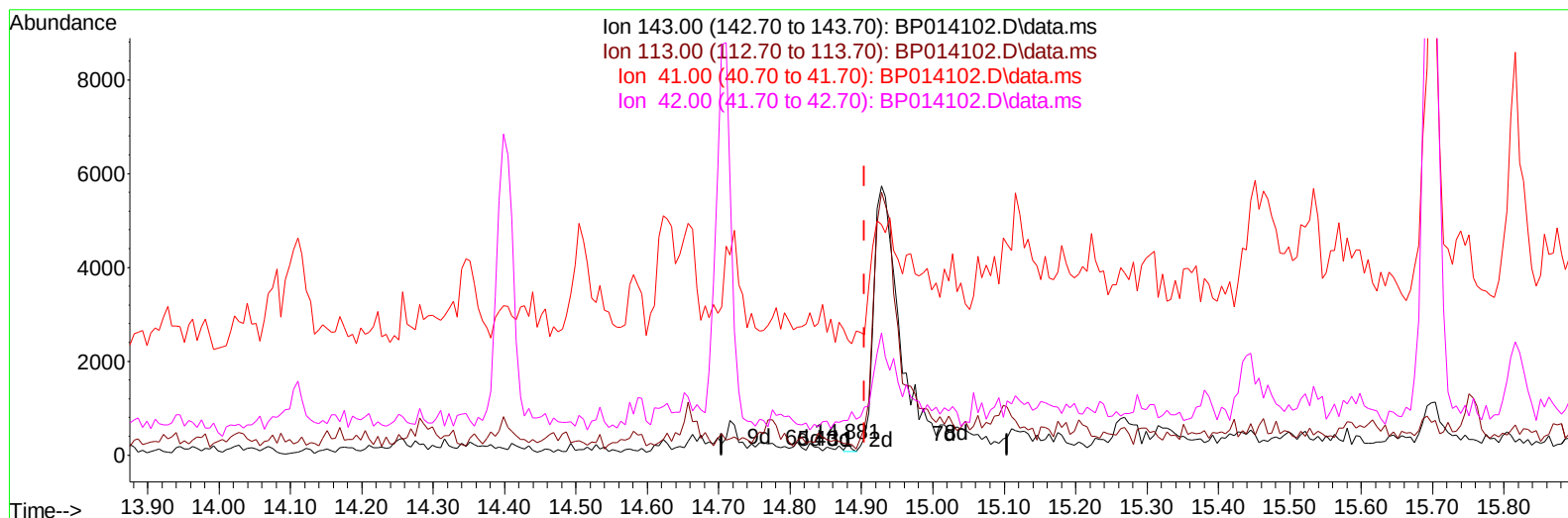
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP031623\
 Data File : BP014102.D
 Acq On : 16 Mar 2023 14:45
 Operator : CG/JU
 Sample : 01884-03
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 CB908

Manual IntegrationsAPPROVED

Quant Time: Mar 17 00:59:53 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP030723.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Mar 17 00:57:25 2023
 Response via : Initial Calibration

Reviewed By :Christian Giraldo 03/17/2023
 Supervised By :Jagrut Upadhyay 03/17/2023



TIC: BP014102.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.881min (-0.023) 0.01 ng/ul

response 74

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	103.10	98.79
41.00	40.20	991.90#
42.00	28.50	298.79#

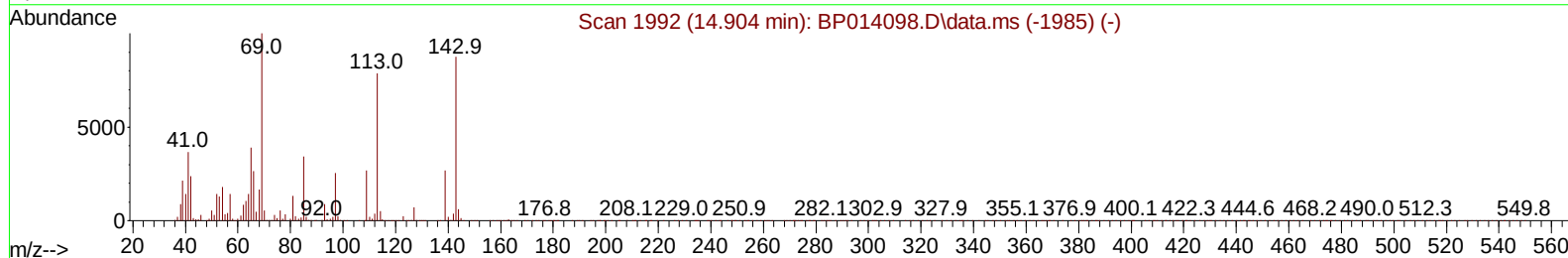
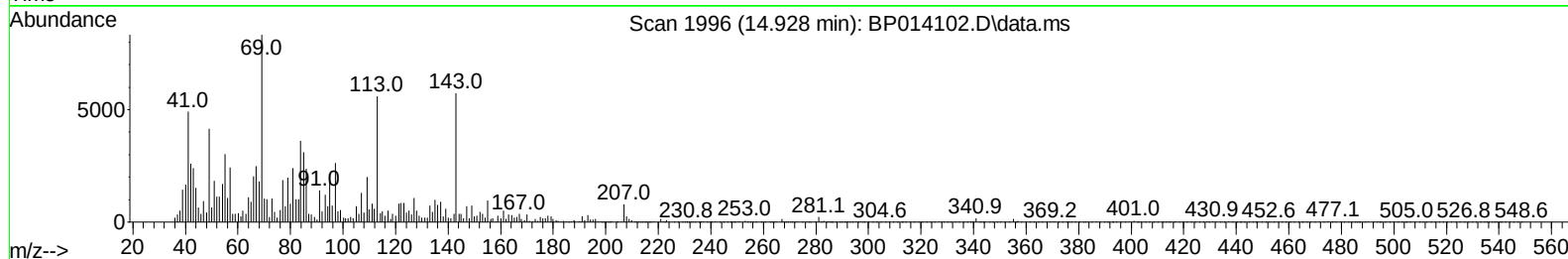
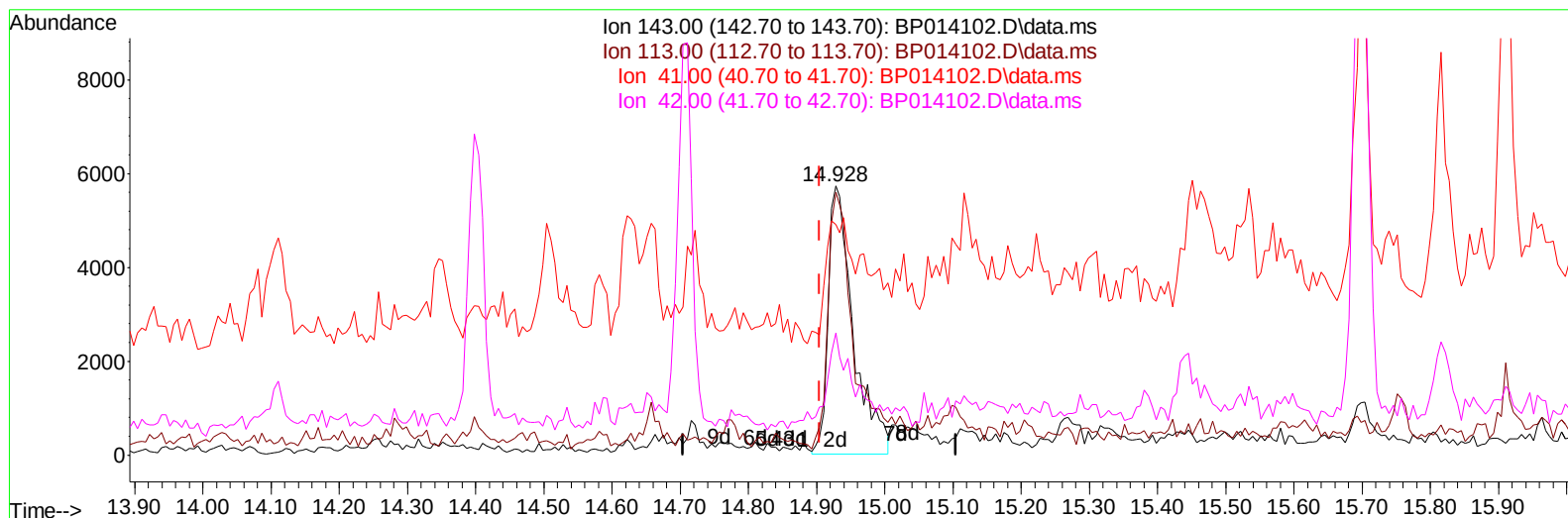
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(54) 4-Nitrophenol-d4 (S)

14.928min (+ 0.024) 2.80 ng/ul m

response 14619

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	103.10	97.84
41.00	40.20	85.76#
42.00	28.50	45.31#

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.063	152	152688	20.000	ng/ul	0.00
20) Naphthalene-d8	10.887	136	518187	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.710	164	349148	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.463	188	939345	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.545	240	1093049	20.000	ng/ul	# 0.00
88) Perylene-d12	24.080	264	1129278	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.422	96	14400	3.563	ng/uL	0.00
4) Pyridine-d5	3.864	84	112124	10.021	ng/ul	0.00
7) Phenol-d5	7.216	99	49477	3.672	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.381	67	188449	23.981	ng/ul	0.00
11) 2-Chlorophenol-d4	7.593	132	114710	11.174	ng/ul	0.00
15) 4-Methylphenol-d8	8.775	113	109599	9.958	ng/ul	0.00
21) Nitrobenzene-d5	9.240	128	111580	26.987	ng/ul	0.00
24) 2-Nitrophenol-d4	9.963	143	64399	13.012	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.504	165	123158	13.225	ng/ul	0.00
31) 4-Chloroaniline-d4	11.028	131	201539	16.353	ng/ul	0.00
46) Dimethylphthalate-d6	14.110	166	728920	24.600	ng/ul	0.00
49) Acenaphthylene-d8	14.398	160	716637	22.830	ng/ul	0.00
54) 4-Nitrophenol-d4	14.928	143	14619m	2.800	ng/ul	0.02
60) Fluorene-d10	15.698	176	726736	28.939	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.816	200	77471	10.912	ng/ul	0.00
73) Anthracene-d10	17.563	188	1219893	26.752	ng/ul	0.00
81) Pyrene-d10	19.786	212	1551042	25.728	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.916	264	1498067	24.939	ng/ul	0.00
Target Compounds						
						Qvalue
52) Acenaphthene	14.769	153	61612	2.675	ng/ul	98
61) Fluorene	15.757	166	60443	2.176	ng/ul	96

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

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