

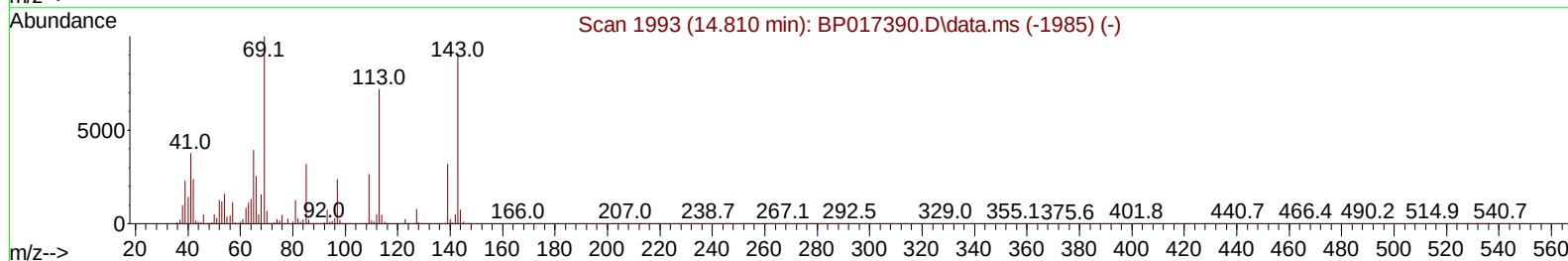
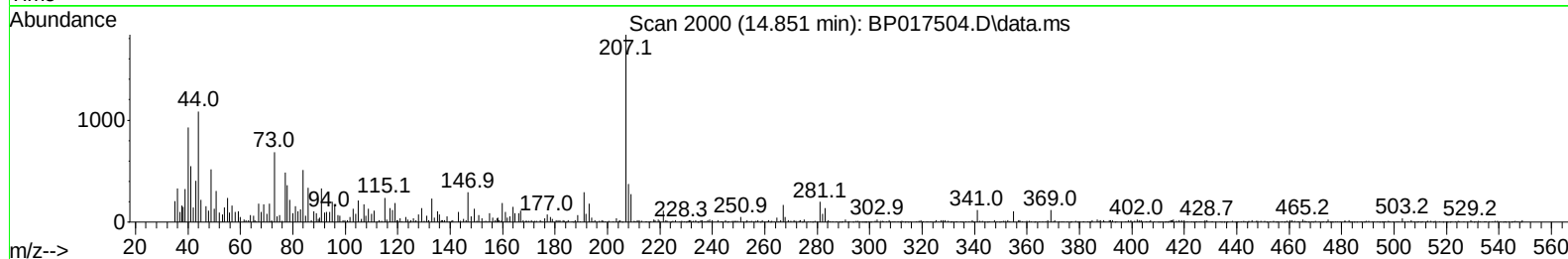
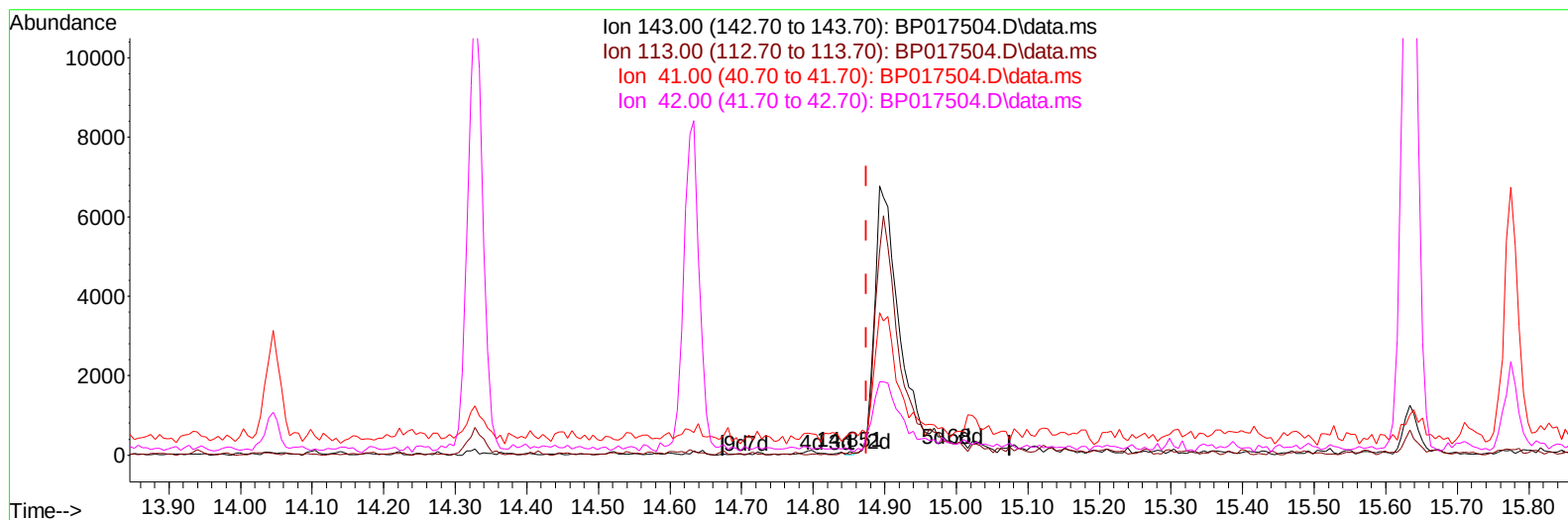
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100223\
 Data File : BP017504.D
 Acq On : 03 Oct 2023 07:01
 Operator : MA/JU
 Sample : 04588-06
 Misc :
 ALS Vial : 28 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BBJE3

Manual IntegrationsAPPROVED

Quant Time: Oct 03 07:37:49 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092023.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Oct 03 07:02:34 2023
 Response via : Initial Calibration

Reviewed By :Yogesh Patel 10/03/2023
 Supervised By :mohammad ahmed 10/05/2023



TIC: BP017504.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.851min (-0.023) 0.01 ng/u1

response 39

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	74.90	20.62#
41.00	36.70	567.01#
42.00	26.40	145.36#

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Quant Time: Oct 03 07:38:28 2023
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.934	152	140844	20.000	ng/ul	-0.01
20) Naphthalene-d8	10.775	136	570776	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.634	164	333055	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.422	188	684919	20.000	ng/ul	0.00
79) Chrysene-d12	21.557	240	667849	20.000	ng/ul	0.00
88) Perylene-d12	24.145	264	743779	20.000	ng/ul	0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.287	96	18230	5.316	ng/uL	0.00
4) Pyridine-d5	3.722	84	133367	13.912	ng/ul	0.00
7) Phenol-d5	7.093	99	88069	7.613	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.275	67	290856	41.665	ng/ul	-0.01
11) 2-Chlorophenol-d4	7.457	132	283322	30.881	ng/ul	0.00
15) 4-Methylphenol-d8	8.657	113	170029	18.906	ng/ul	-0.01
21) Nitrobenzene-d5	9.146	128	184235	44.648	ng/ul	0.00
24) 2-Nitrophenol-d4	9.857	143	187374	41.820	ng/ul	-0.01
28) 2,4-Dichlorophenol-d3	10.393	165	317130	37.464	ng/ul	0.00
31) 4-Chloroaniline-d4	10.946	131	394826	32.199	ng/ul	0.00
46) Dimethylphthalate-d6	14.045	166	1048247	43.935	ng/ul	0.00
49) Acenaphthylene-d8	14.328	160	1176569	42.877	ng/ul	-0.01
54) 4-Nitrophenol-d4	14.892	143	17254m	4.151	ng/ul	0.02
60) Fluorene-d10	15.634	176	875652	42.631	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.775	200	91528	23.710	ng/ul	0.00
73) Anthracene-d10	17.522	188	1406891	46.315	ng/ul	0.00
81) Pyrene-d10	19.774	212	1777993	49.231	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.974	264	1765701	49.306	ng/ul	0.02

Target Compounds Qvalue

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

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