

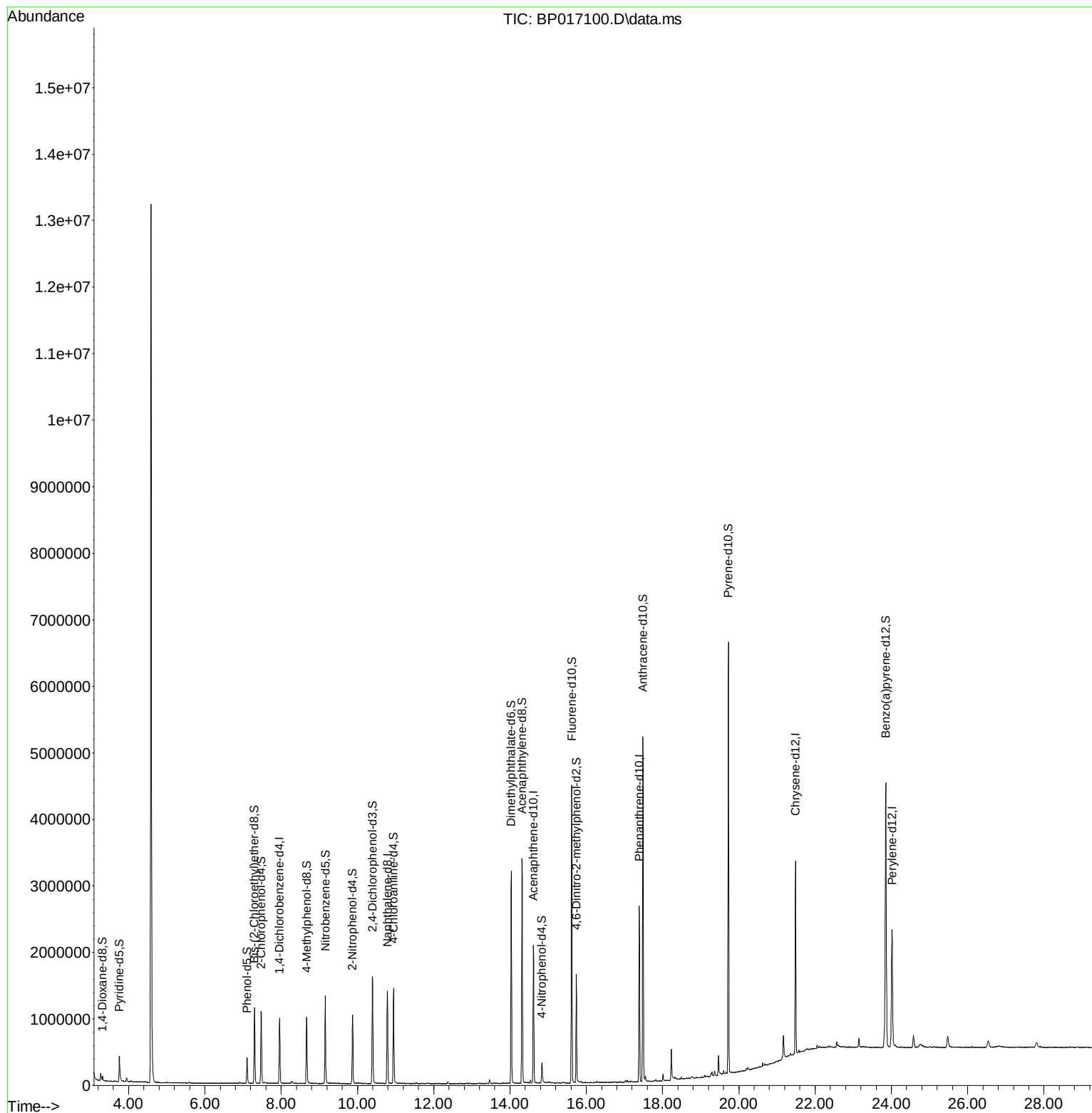
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP091023\
Data File : BP017100.D
Acq On : 11 Sep 2023 12:31
Operator : MA/JU
Sample : 04262-06
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BBHM8

Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 09/12/2023
Supervised By :mohammad ahmed 09/12/2023

Quant Time: Sep 12 02:41:21 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP091023.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Mon Sep 11 06:13:29 2023
Response via : Initial Calibration



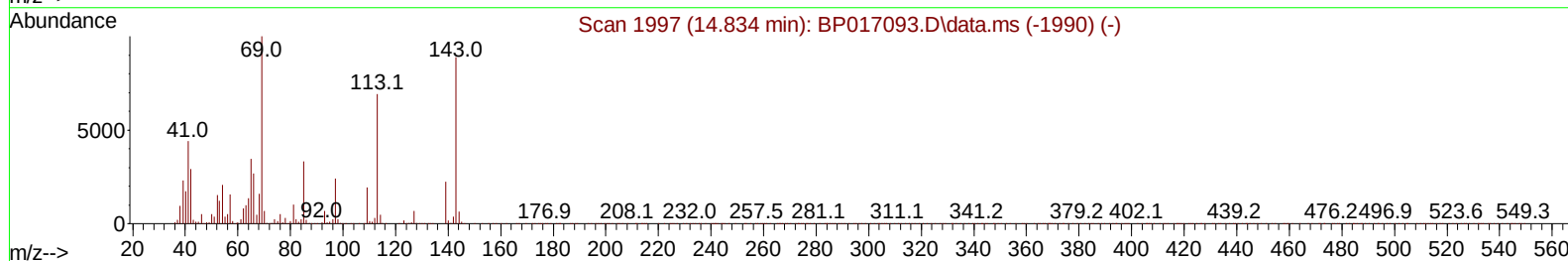
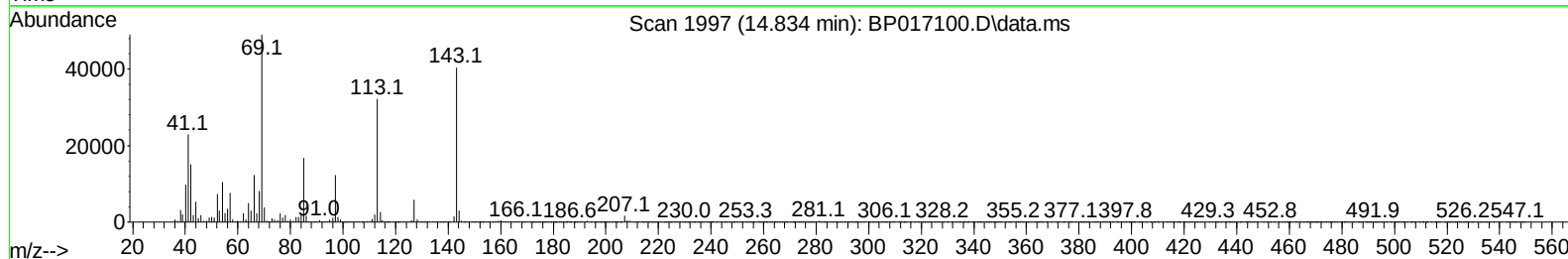
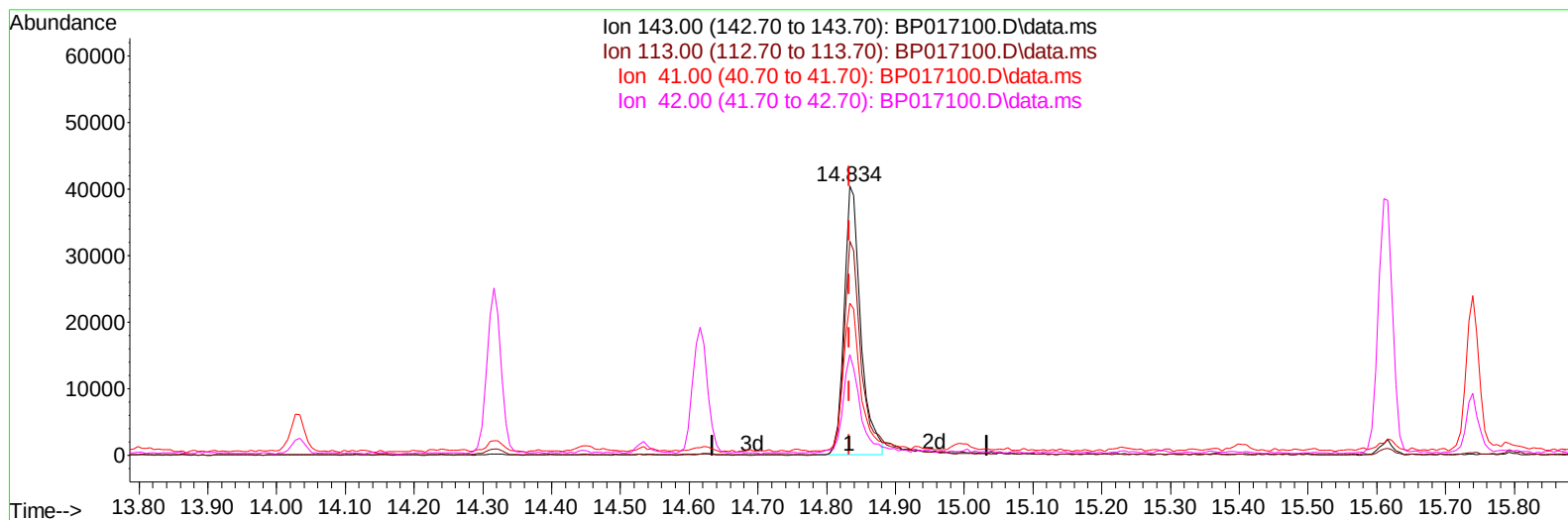
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TIC: BP017100.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.834min (+ 0.000) 7.24 ng/ul

response 67564

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	76.70	79.48
41.00	51.90	56.48
42.00	35.60	37.44

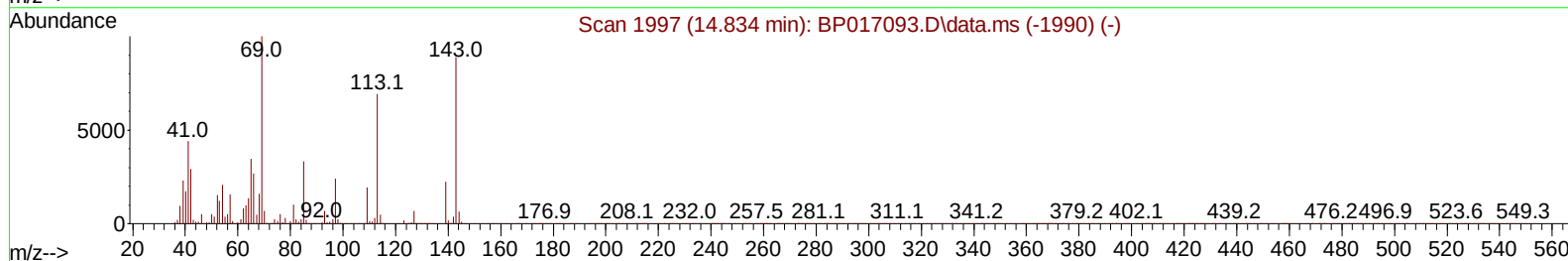
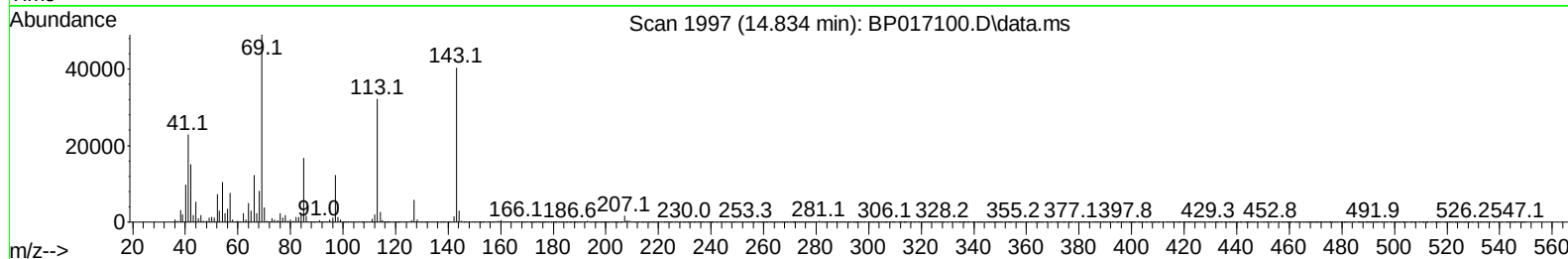
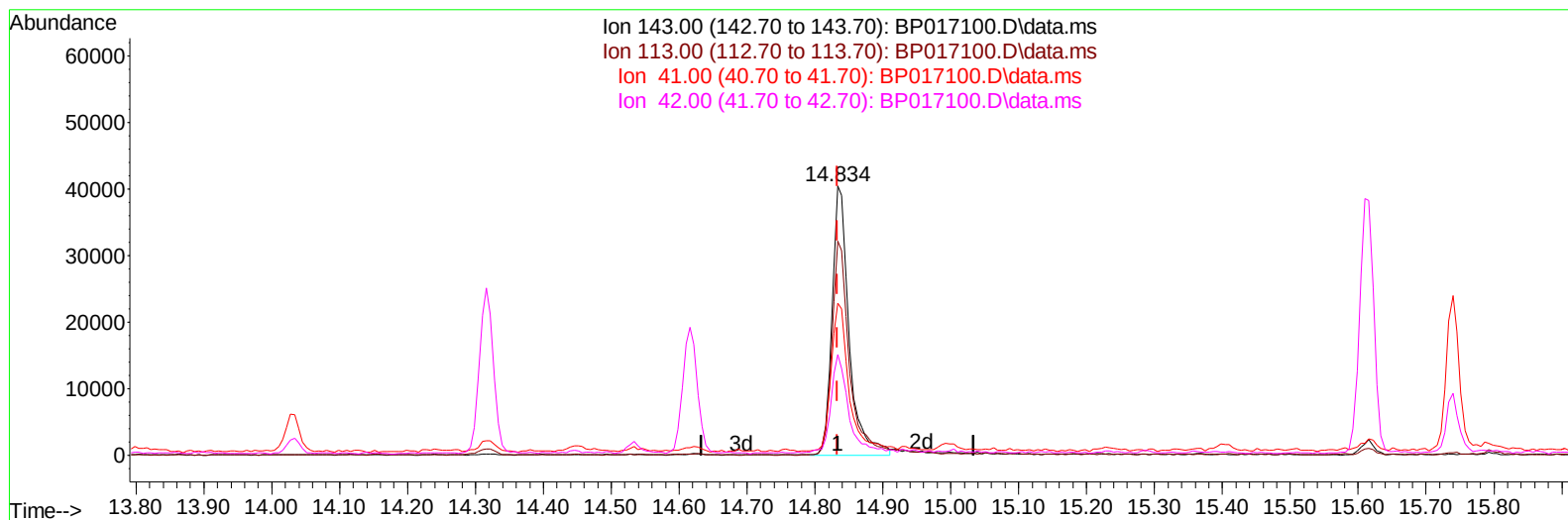
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TIC: BP017100.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.834min (+ 0.000) 7.53 ng/ul m

response 70269

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	76.70	79.48
41.00	51.90	56.48
42.00	35.60	37.44

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.958	152	235779	20.000	ng/ul	0.00
20) Naphthalene-d8	10.787	136	1059828	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.616	164	635338	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.386	188	1368776	20.000	ng/ul	0.00
79) Chrysene-d12	21.486	240	1166438	20.000	ng/ul	0.00
88) Perylene-d12	24.015	264	1207417	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.317	96	35567	5.758	ng/uL	0.00
4) Pyridine-d5	3.752	84	217495	11.676	ng/ul	0.00
7) Phenol-d5	7.111	99	205413	9.584	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.299	67	492533	34.264	ng/ul	0.00
11) 2-Chlorophenol-d4	7.481	132	443524	28.632	ng/ul	0.00
15) 4-Methylphenol-d8	8.669	113	354944	20.935	ng/ul	0.00
21) Nitrobenzene-d5	9.157	128	262589	34.801	ng/ul	0.00
24) 2-Nitrophenol-d4	9.869	143	270817	34.776	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.393	165	508236	34.418	ng/ul	0.00
31) 4-Chloroaniline-d4	10.946	131	670514	28.317	ng/ul	0.00
46) Dimethylphthalate-d6	14.034	166	1892246	41.857	ng/ul	0.00
49) Acenaphthylene-d8	14.316	160	1995447	38.764	ng/ul	0.00
54) 4-Nitrophenol-d4	14.834	143	70269m	7.534	ng/ul	0.00
60) Fluorene-d10	15.616	176	1576697	41.099	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.739	200	237214	30.709	ng/ul	0.00
73) Anthracene-d10	17.486	188	2540432	43.300	ng/ul	0.00
81) Pyrene-d10	19.727	212	2979644	48.507	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.851	264	2612093	45.064	ng/ul	0.00

Target Compounds Qvalue

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