

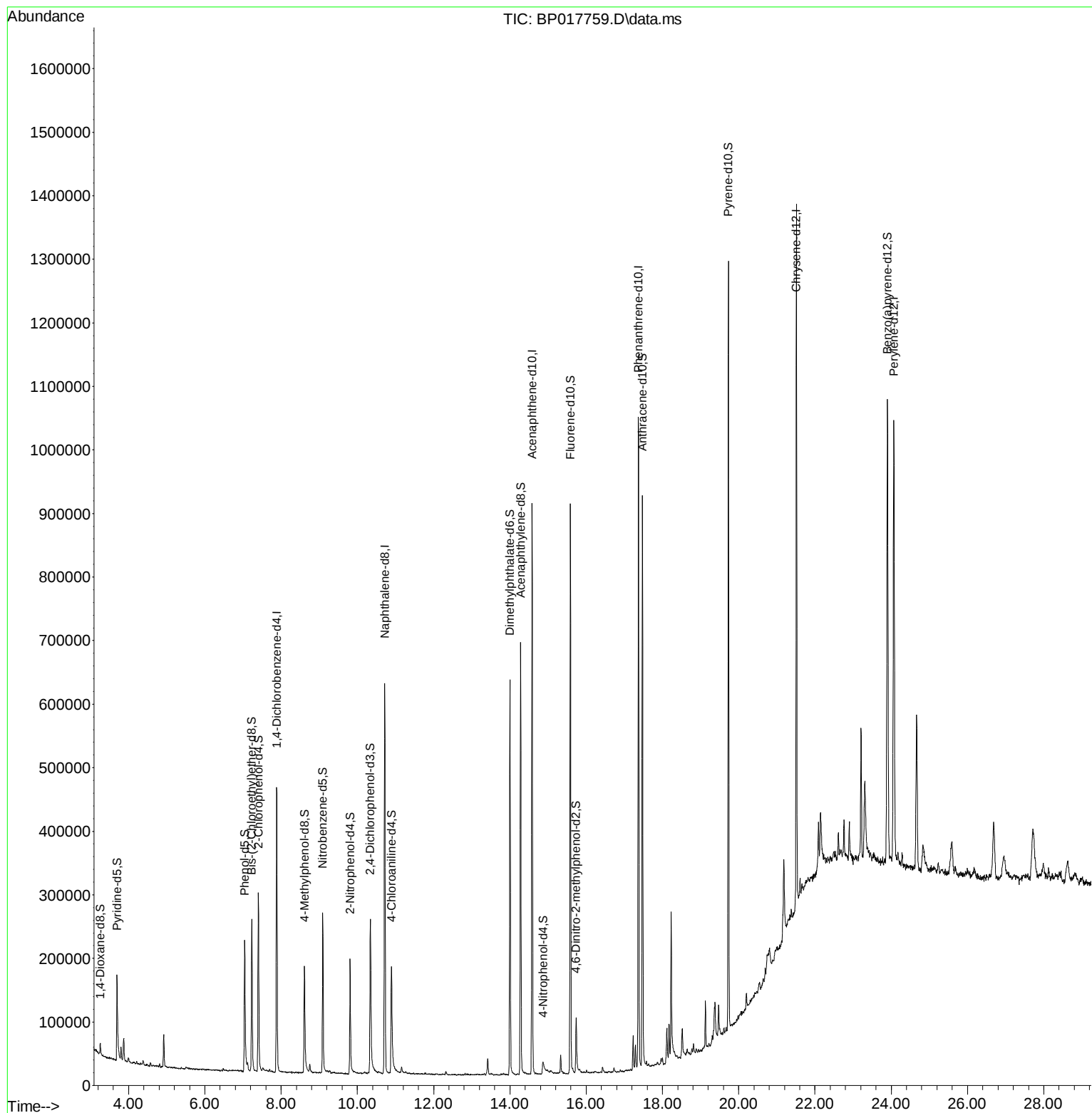
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102323\
Data File : BP017759.D
Acq On : 23 Oct 2023 20:35
Operator : MA/JU
Sample : 04926-08
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
GCCZ3

Manual IntegrationsAPPROVED

Quant Time: Oct 24 06:03:11 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP101823.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Thu Oct 19 01:32:15 2023
Response via : Initial Calibration

Reviewed By :Yogesh Patel 10/24/2023
Supervised By :mohammad ahmed 10/25/2023



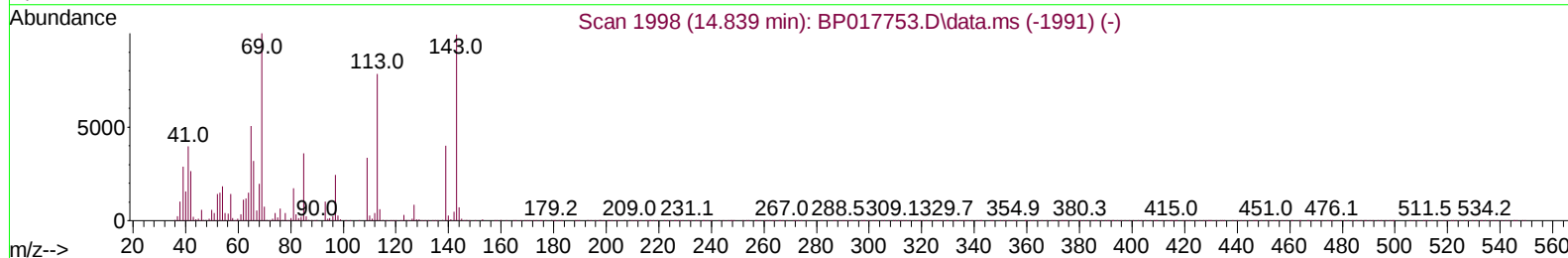
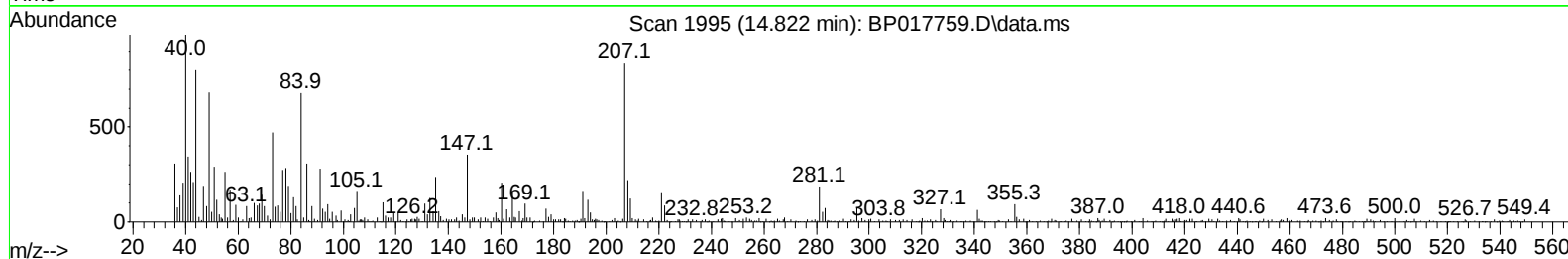
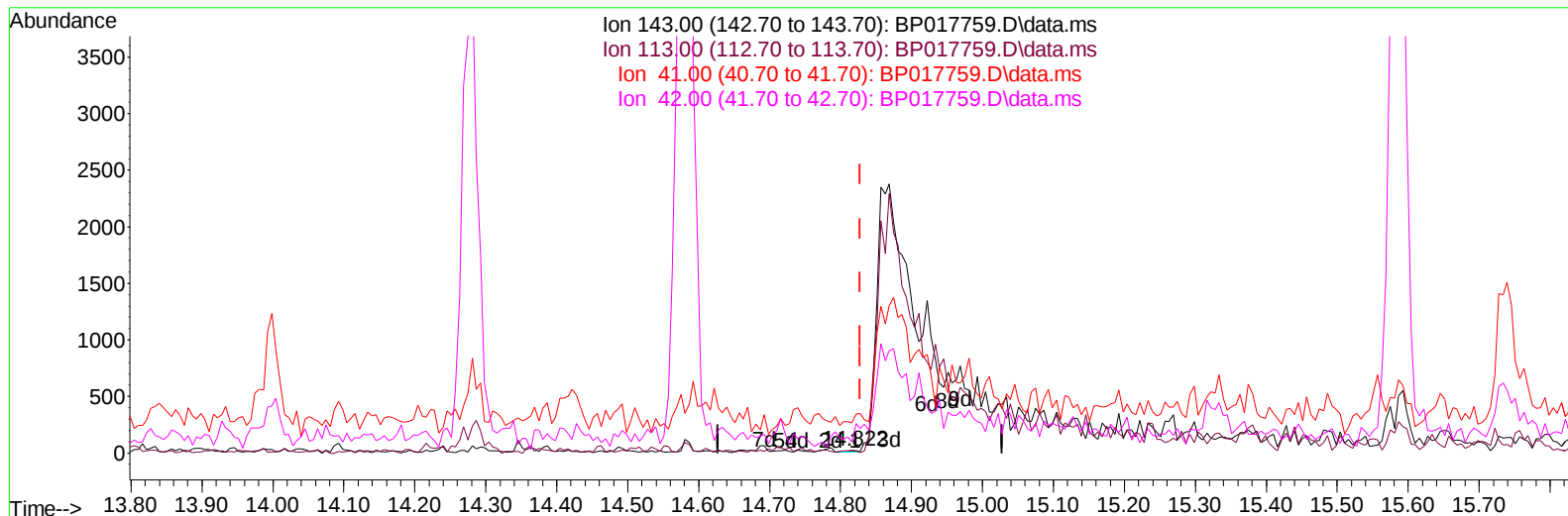
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TIC: BP017759.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.822min (-0.006) 0.00 ng/u1

response

14

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	81.70	100.00#
41.00	39.20	1572.73#
42.00	25.60	1195.45#

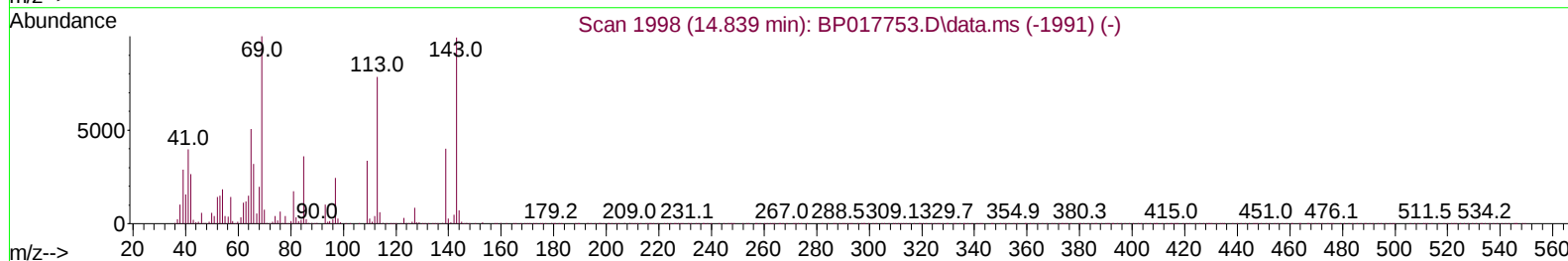
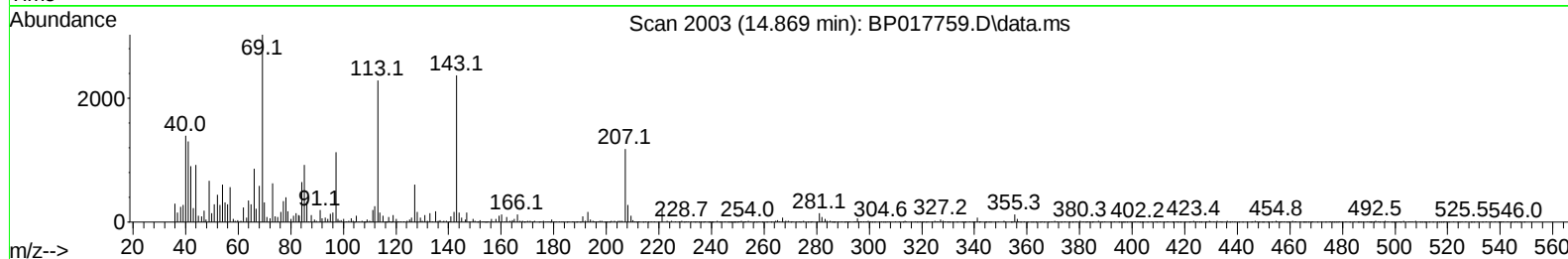
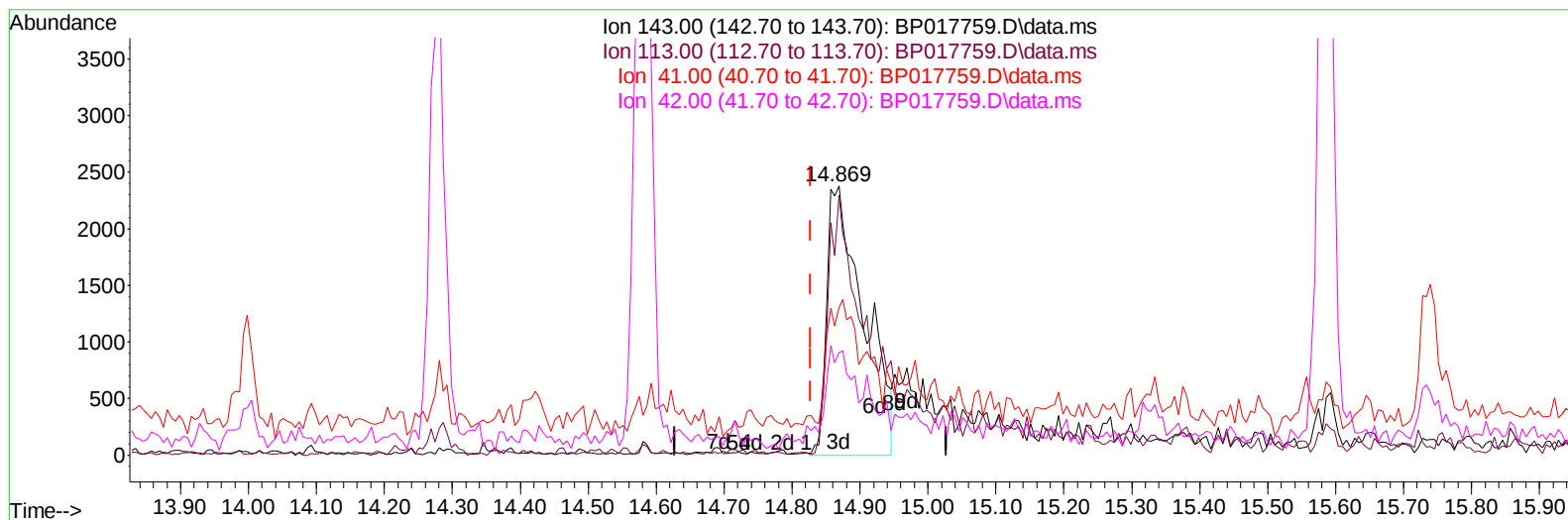
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TIC: BP017759.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.869min (+ 0.041) 2.65 ng/ul m

response 8960

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	81.70	96.47
41.00	39.20	54.86#
42.00	25.60	37.99#

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.887	152	117485	20.000	ng/ul	0.00
20) Naphthalene-d8	10.716	136	473129	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.581	164	282814	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.369	188	599722	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.516	240	478946	20.000	ng/ul	0.00
88) Perylene-d12	24.068	264	501473	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.258	96	9396	2.975	ng/uL	0.00
4) Pyridine-d5	3.693	84	83754	10.129	ng/ul	0.00
7) Phenol-d5	7.046	99	122559	12.536	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.228	67	104657	17.334	ng/ul	0.00
11) 2-Chlorophenol-d4	7.405	132	120469	15.108	ng/ul	0.00
15) 4-Methylphenol-d8	8.610	113	72022	9.381	ng/ul	0.00
21) Nitrobenzene-d5	9.093	128	55909	14.583	ng/ul	0.00
24) 2-Nitrophenol-d4	9.810	143	54112	12.774	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.340	165	99664	11.994	ng/ul	0.00
31) 4-Chloroaniline-d4	10.898	131	114281	10.509	ng/ul	0.00
46) Dimethylphthalate-d6	13.998	166	392959	16.552	ng/ul	0.00
49) Acenaphthylene-d8	14.281	160	427573	16.149	ng/ul	0.00
54) 4-Nitrophenol-d4	14.869	143	8960m	2.649	ng/ul	0.04
60) Fluorene-d10	15.586	176	333601	16.255	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.734	200	20690	5.238	ng/ul	0.00
73) Anthracene-d10	17.475	188	498520	16.266	ng/ul	0.00
81) Pyrene-d10	19.733	212	597600	19.902	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.898	264	491370	17.593	ng/ul	0.00

Target Compounds Qvalue

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