

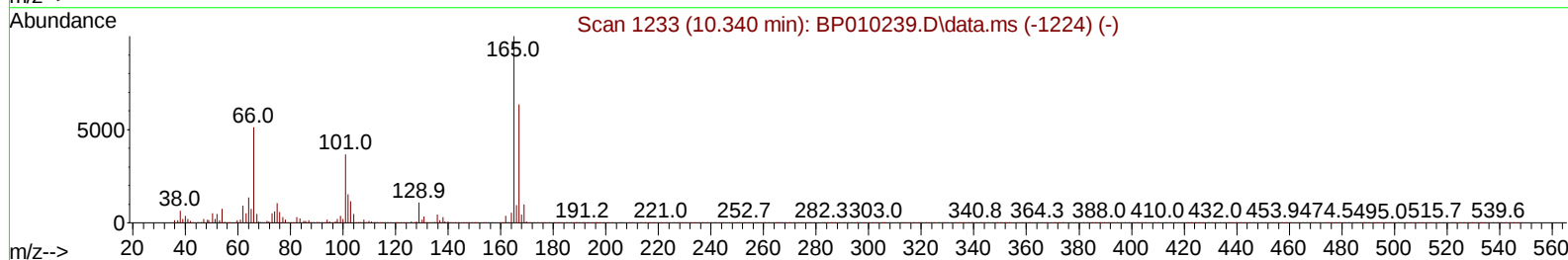
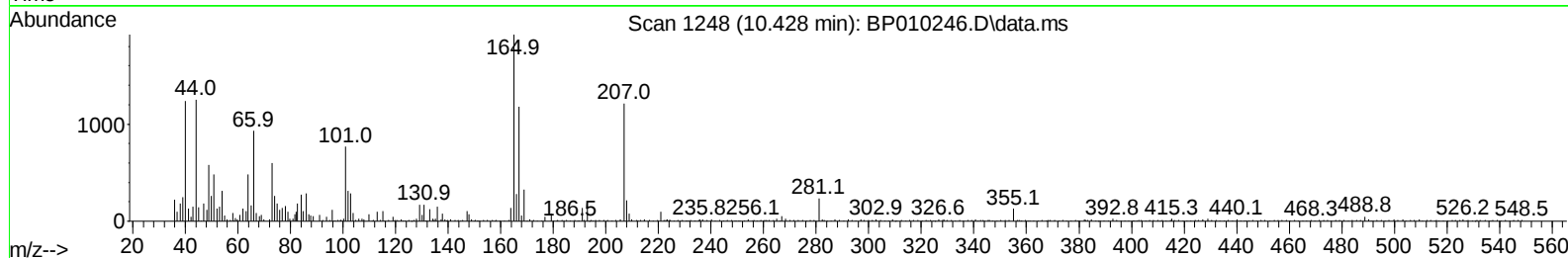
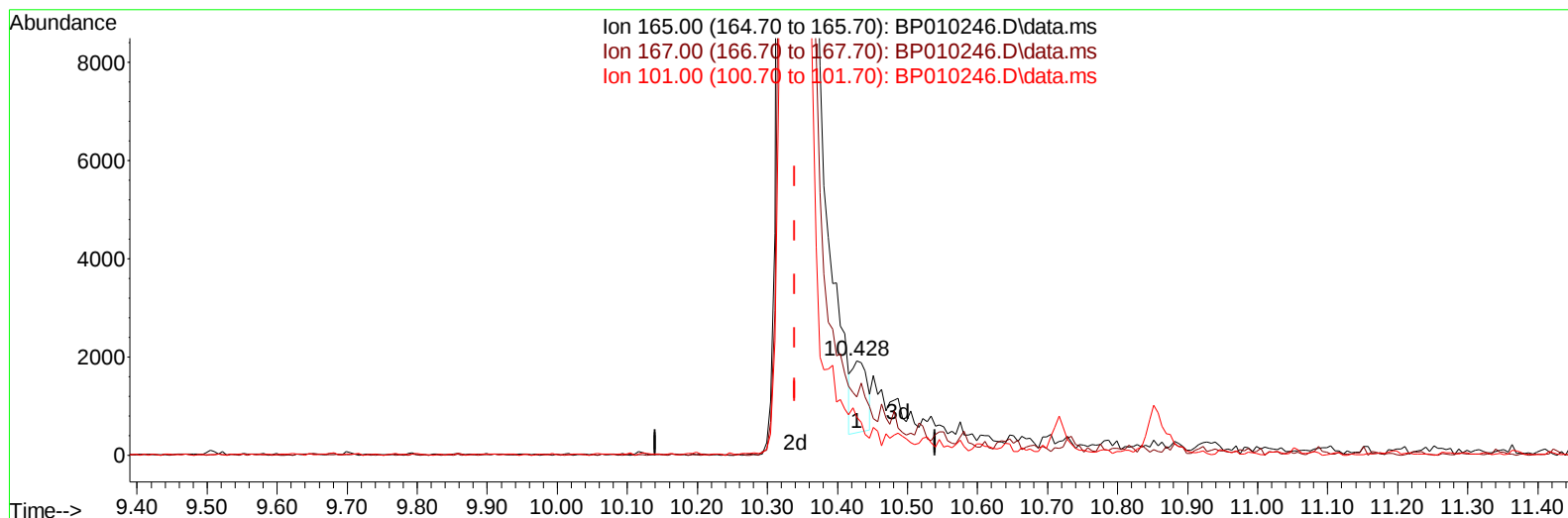
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP050522\
Data File : BP010246.D
Acq On : 05 May 2022 14:36
Operator : CG/JU
Sample : N2668-17
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EXL68

Manual IntegrationsAPPROVED

Quant Time: May 06 01:33:10 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP042122.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri May 06 01:05:07 2022
Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 05/10/2022
Supervised By :mohammad ahmed 05/10/2022



TIC: BP010246.D\data.ms

(28) 2,4-Dichlorophenol-d3 (S)

10.428min (+ 0.088) 0.20 ng/u1

response 2172

Ion	Exp%	Act%
165.00	100.00	100.00
167.00	57.70	61.45
101.00	49.80	39.85
0.00	0.00	0.00

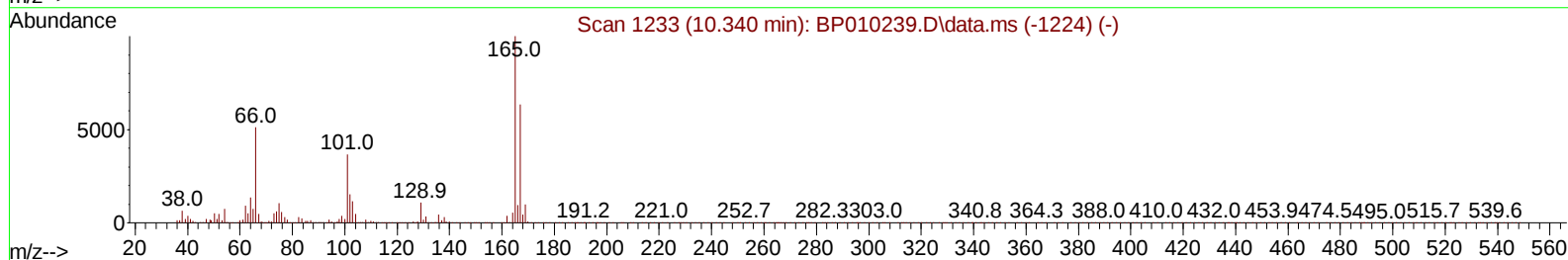
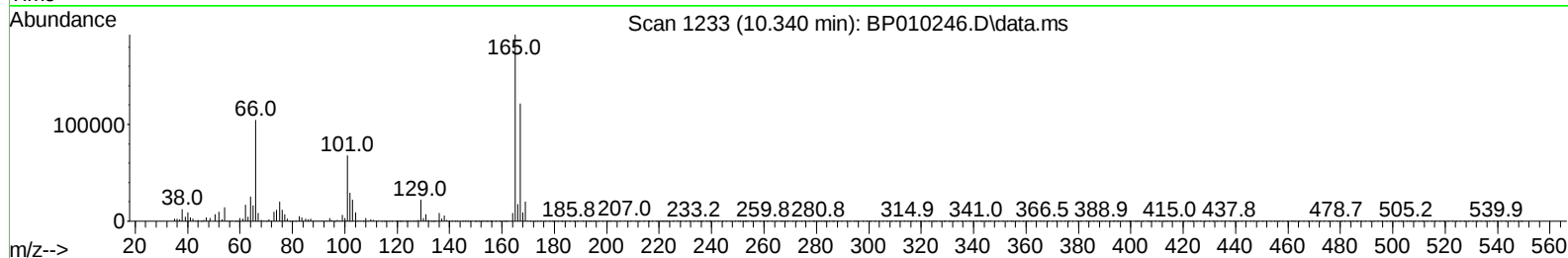
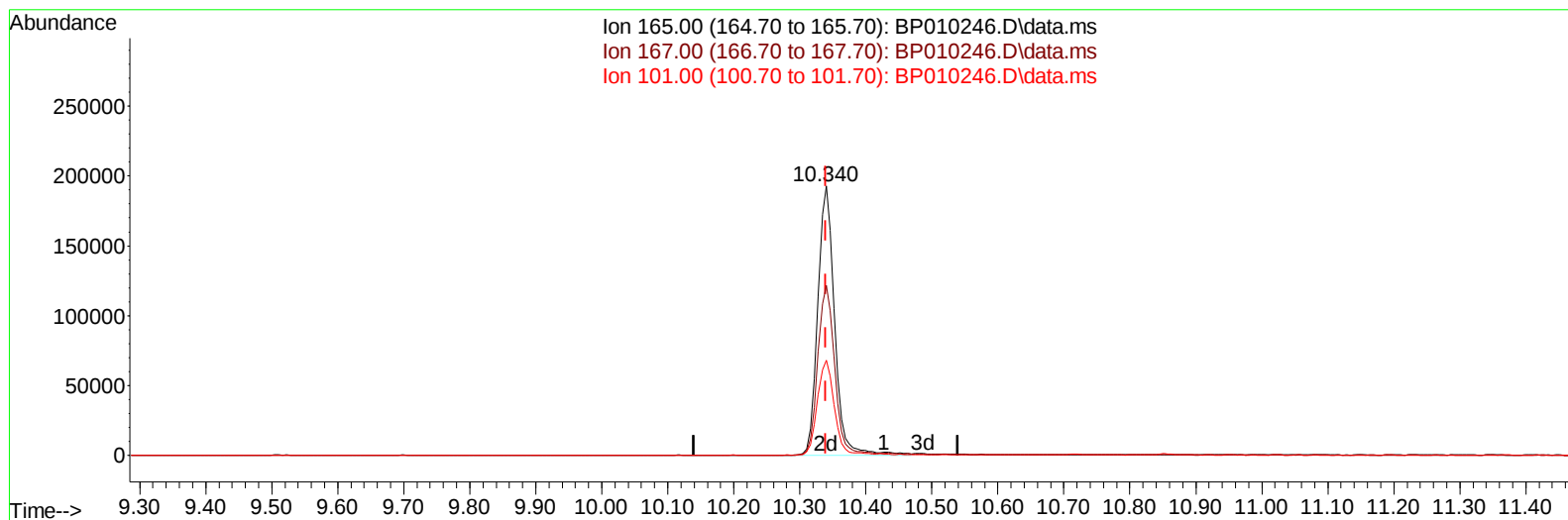
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Instrument :
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TIC: BP010246.D\data.ms

(28) 2,4-Dichlorophenol-d3 (S)

10.340min (+ 0.000) 31.31 ng/ul m

response 340230

Ion	Exp%	Act%
165.00	100.00	100.00
167.00	57.70	63.18
101.00	49.80	35.18#
0.00	0.00	0.00

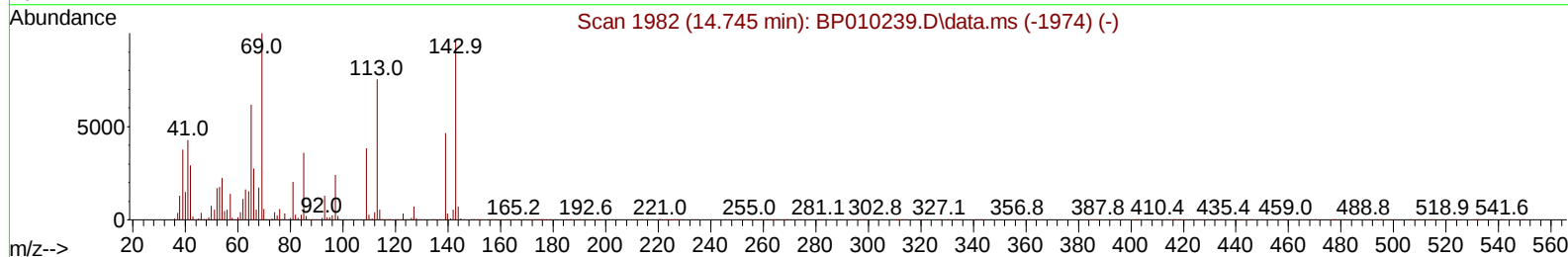
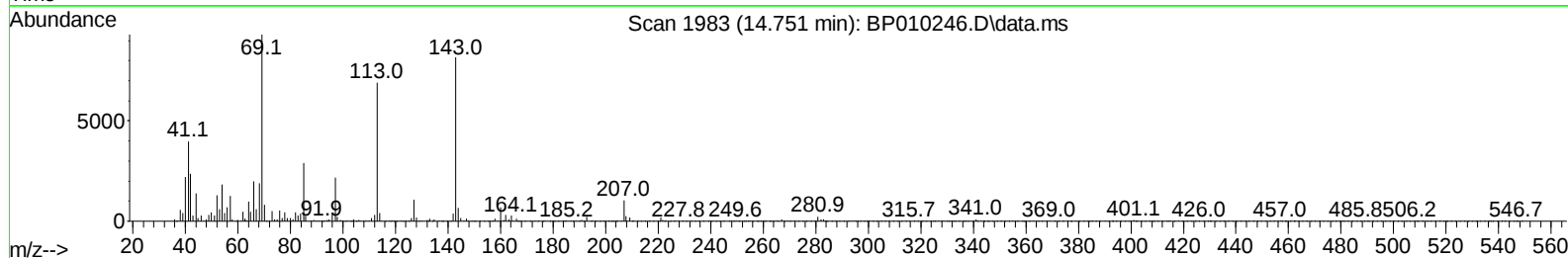
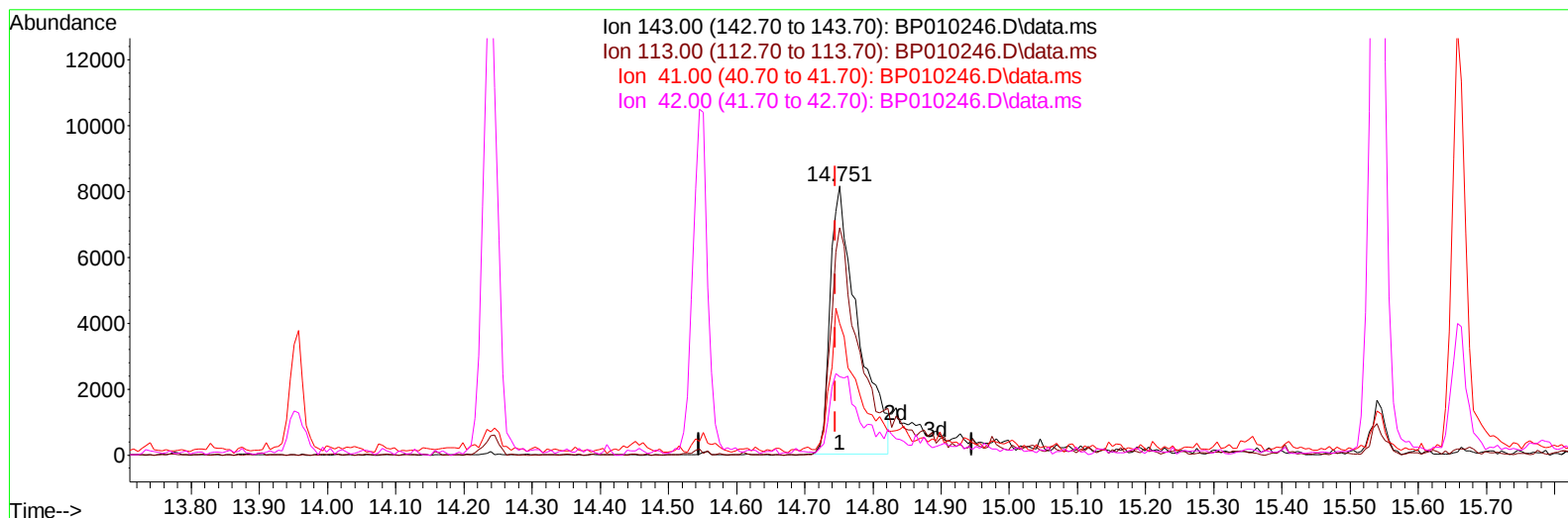
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Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EXL68

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TIC: BP010246.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.751min (+ 0.006) 3.72 ng/u1

response 23036

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	132.70	84.58#
41.00	23.20	48.85#
42.00	18.00	29.06#

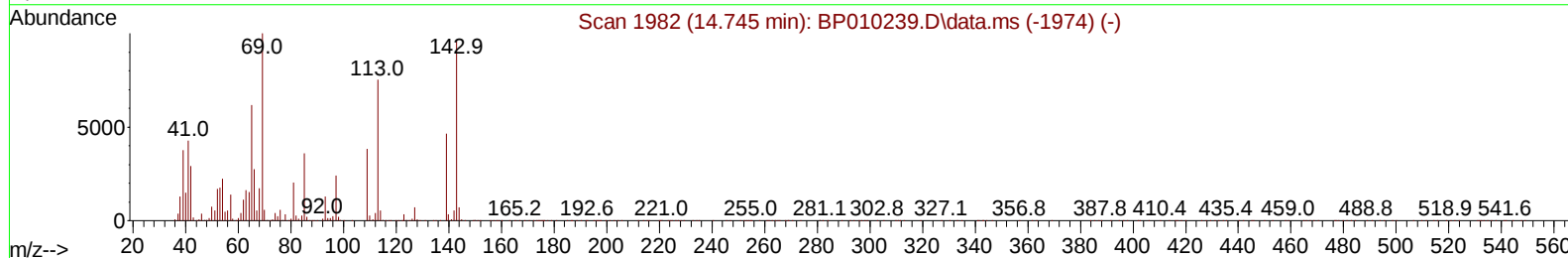
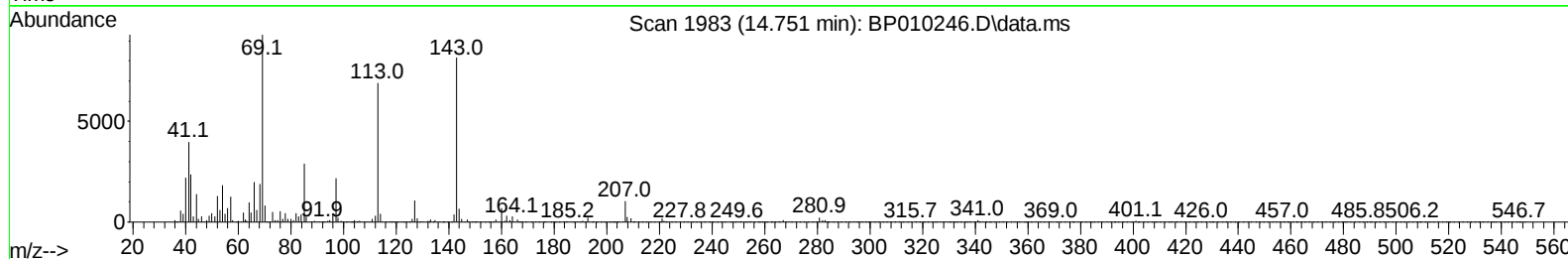
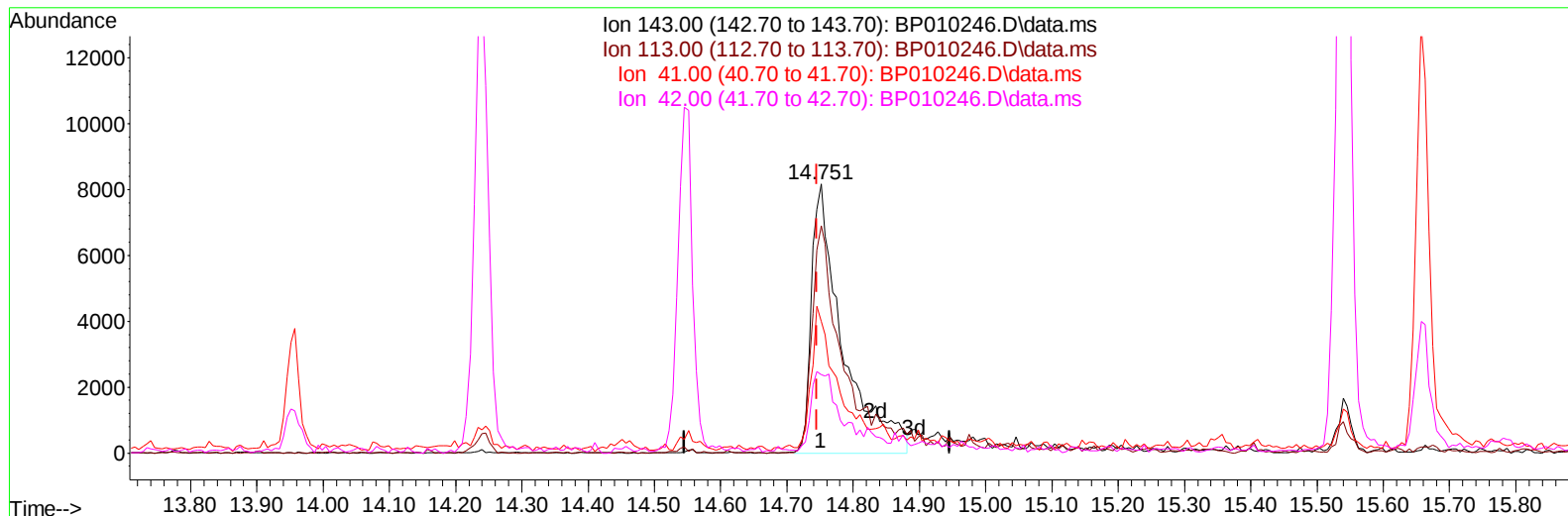
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Sample : N2668-17
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EXL68

Manual IntegrationsAPPROVED

Quant Time: May 06 01:33:10 2022
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TIC: BP010246.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.751min (+ 0.006) 4.29 ng/ul m

response 26573

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	132.70	84.58#
41.00	23.20	48.85#
42.00	18.00	29.06#

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 Misc :
 ALS Vial : 9 Sample Multiplier: 1

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.916	152	142943	20.000	ng/ul	0.00
20) Naphthalene-d8	10.716	136	642603	20.000	ng/ul	# 0.00
38) Acenaphthene-d10	14.545	164	432909	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.298	188	883815	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.380	240	599081	20.000	ng/ul	0.00
88) Perylene-d12	23.821	264	534260	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.358	96	17090	5.214	ng/uL	0.00
4) Pyridine-d5	3.781	84	80474	8.727	ng/ul	0.01
7) Phenol-d5	7.081	99	89199	6.860	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.240	67	315423	40.430	ng/ul	0.00
11) 2-Chlorophenol-d4	7.446	132	282704	29.359	ng/ul	0.00
15) 4-Methylphenol-d8	8.622	113	187898	17.442	ng/ul	0.00
21) Nitrobenzene-d5	9.075	128	200737	39.276	ng/ul	0.00
24) 2-Nitrophenol-d4	9.799	143	208786	37.405	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.340	165	340230m	31.310	ng/ul	0.00
31) 4-Chloroaniline-d4	10.852	131	473336	30.380	ng/ul	0.00
46) Dimethylphthalate-d6	13.957	166	1420250	42.175	ng/ul	0.00
49) Acenaphthylene-d8	14.240	160	1580278	38.774	ng/ul	0.00
54) 4-Nitrophenol-d4	14.751	143	26573m	4.286	ng/ul	0.00
60) Fluorene-d10	15.540	176	1238848	42.766	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.657	200	179690	32.154	ng/ul	0.00
73) Anthracene-d10	17.398	188	1908921	44.645	ng/ul	0.00
81) Pyrene-d10	19.628	212	1994188	58.532	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.663	264	1334788	47.656	ng/ul	0.00

Target Compounds Qvalue

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

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