

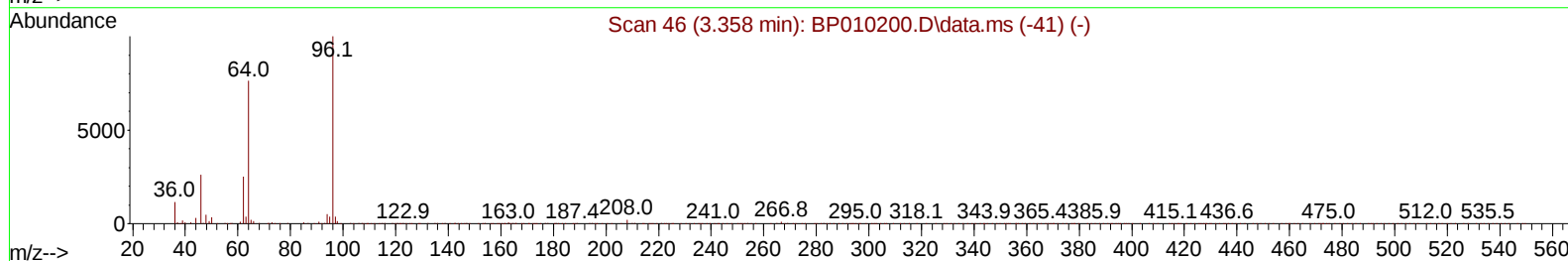
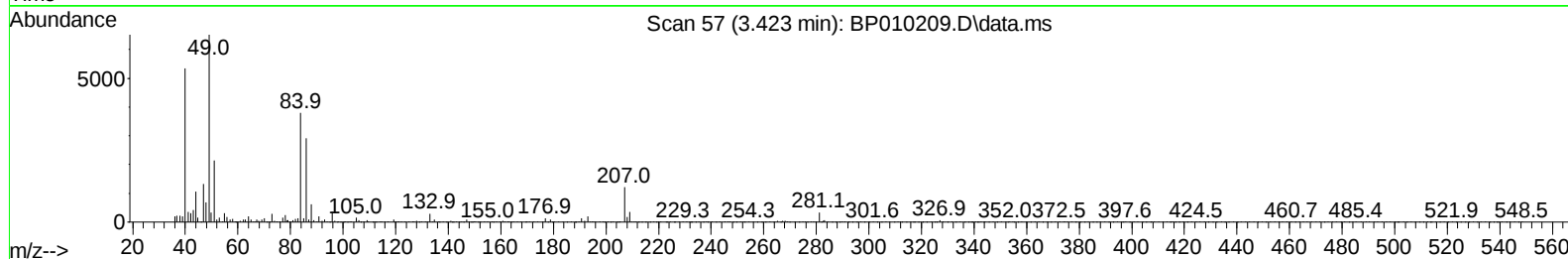
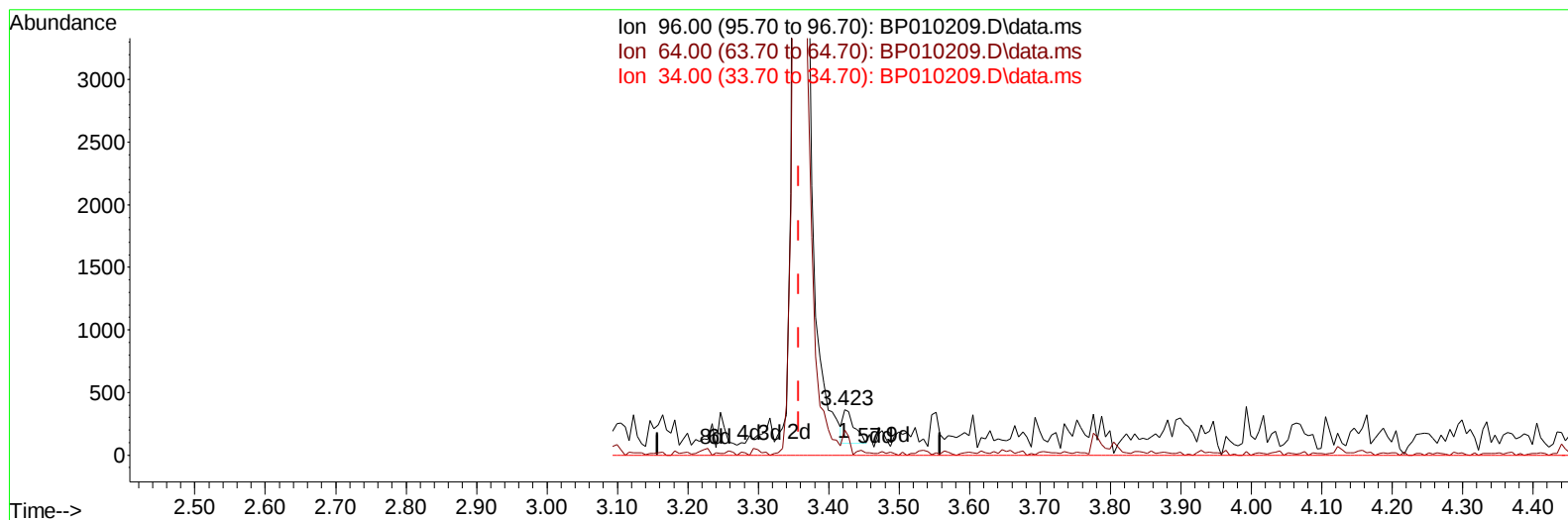
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP050222\
 Data File : BP010209.D
 Acq On : 03 May 2022 16:09
 Operator : CG/JU
 Sample : N2615-17
 Misc :
 ALS Vial : 45 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BH0D4

Manual IntegrationsAPPROVED

Quant Time: May 04 01:10:14 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP042122.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue May 03 00:58:32 2022
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 05/04/2022
 Supervised By :Yogesh Patel 05/05/2022



TIC: BP010209.D\data.ms

(3) 1,4-Dioxane-d8 (S)

3.423min (+ 0.065) 0.11 ng/uL

response 292

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	56.08
34.00	0.00	0.00
0.00	0.00	0.00

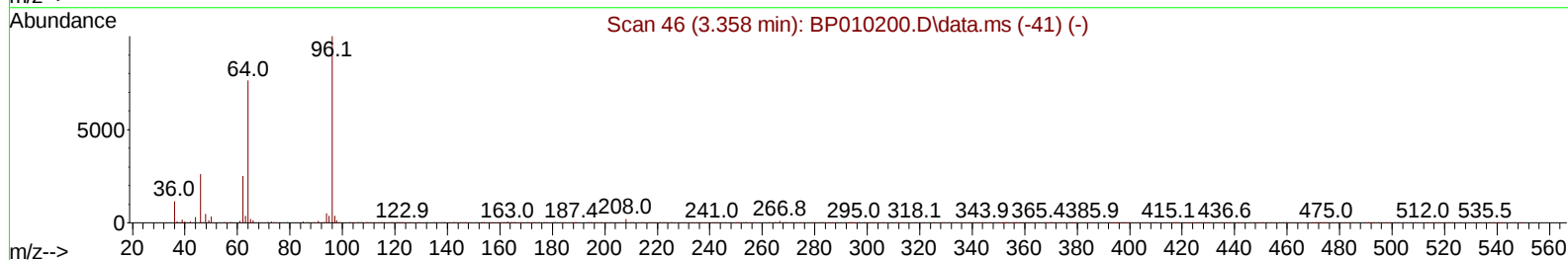
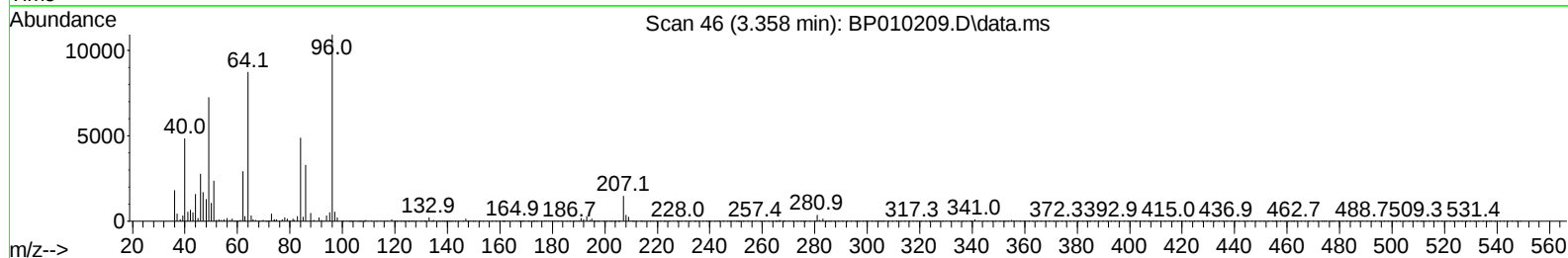
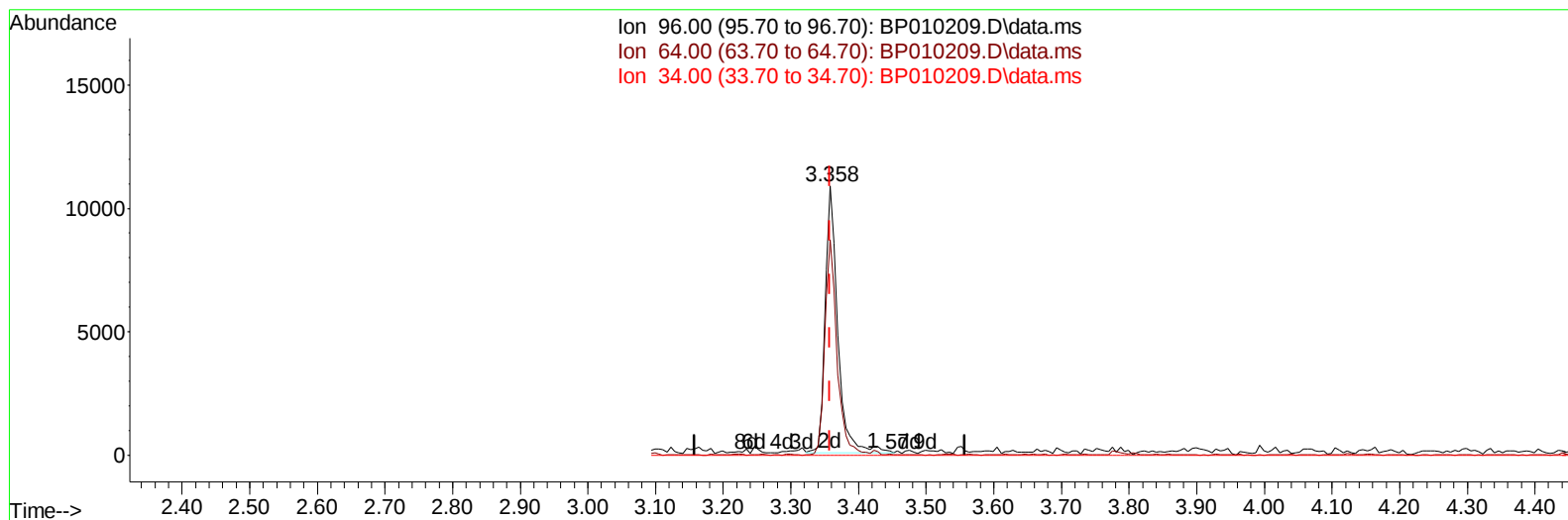
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TIC: BP010209.D\data.ms

(3) 1,4-Dioxane-d8 (S)

3.358min (+ 0.000) 5.39 ng/uL m

response 14171

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	80.05#
34.00	0.00	0.00
0.00	0.00	0.00

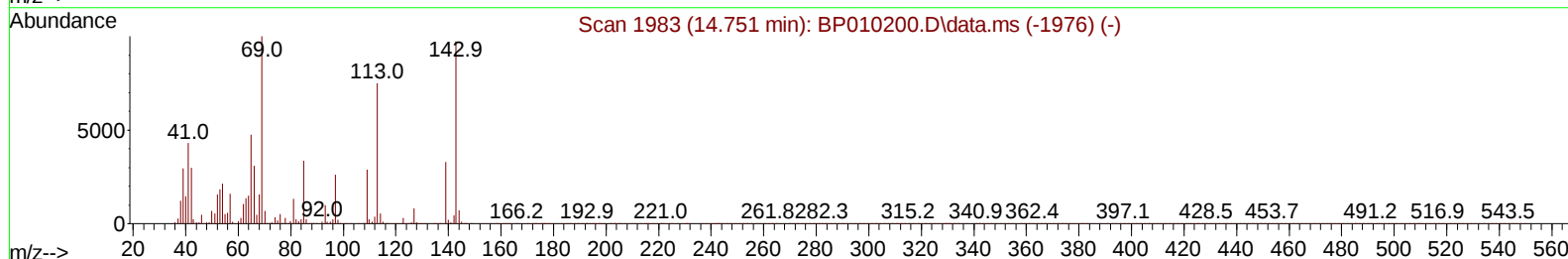
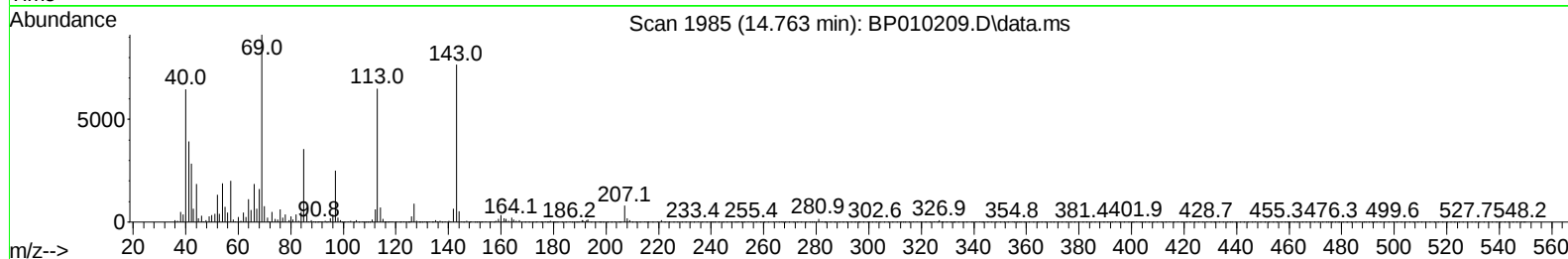
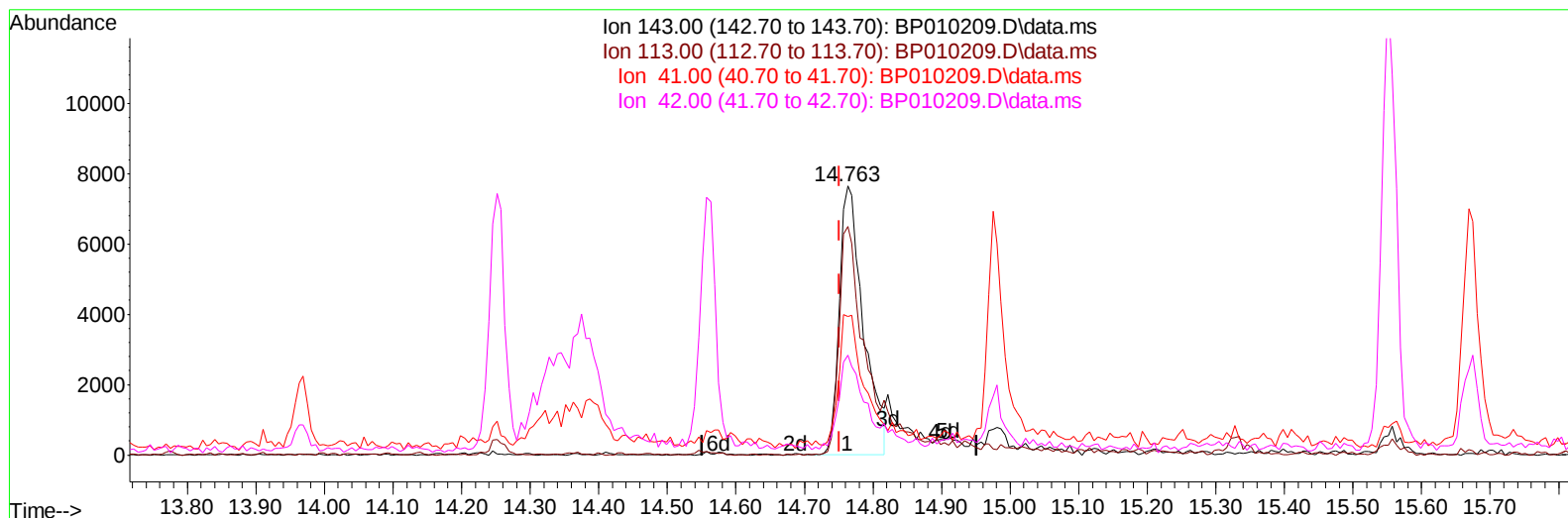
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Misc :
ALS Vial : 45 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BH0D4

Manual IntegrationsAPPROVED

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TIC: BP010209.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.763min (+ 0.012) 4.09 ng/u1

response 18584

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	132.70	84.97#
41.00	23.20	51.51#
42.00	18.00	37.13#

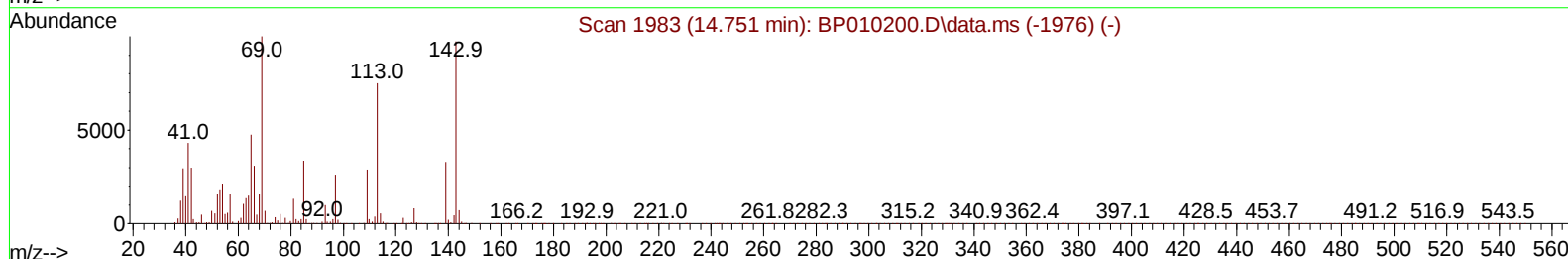
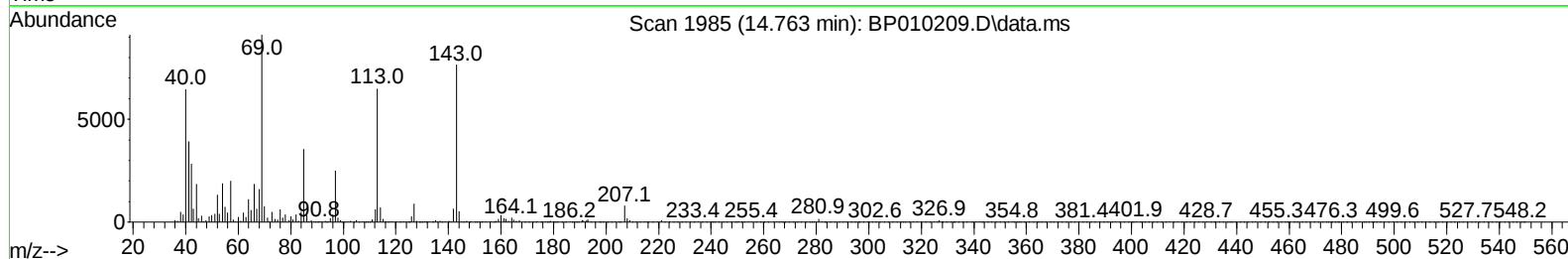
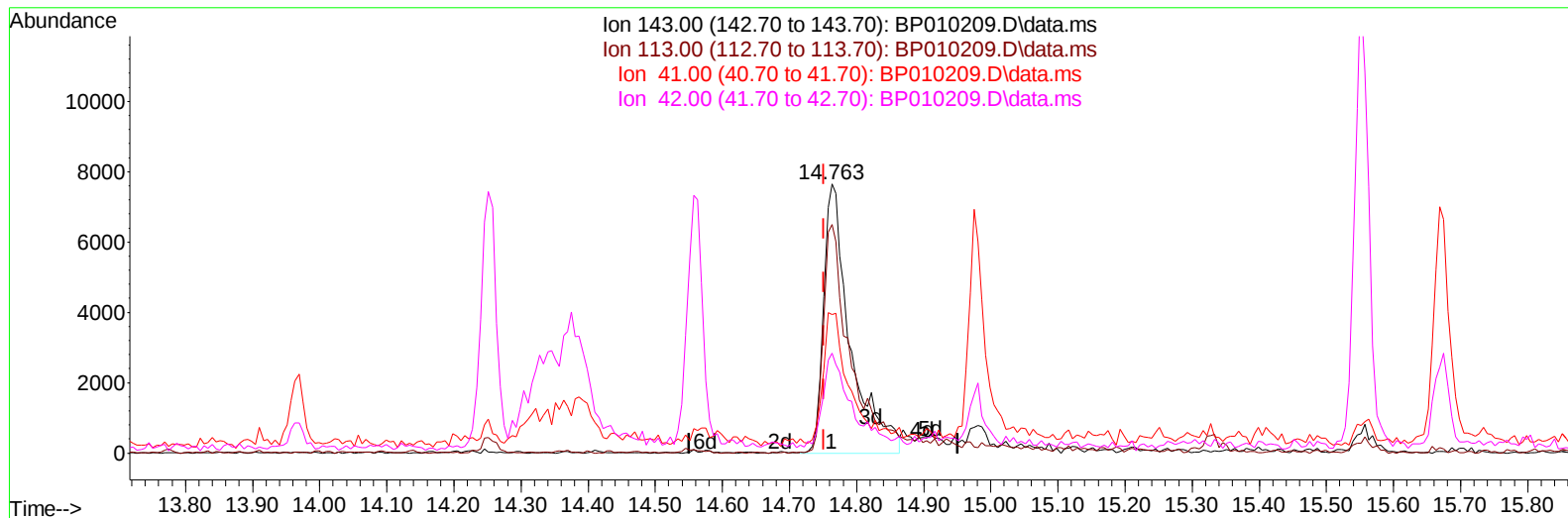
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Acq On : 03 May 2022 16:09
Operator : CG/JU
Sample : N2615-17
Misc :
ALS Vial : 45 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BH0D4

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 05/04/2022
Supervised By :Yogesh Patel 05/05/2022

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TIC: BP010209.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.763min (+ 0.012) 4.66 ng/ul m

response 21171

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	132.70	84.97#
41.00	23.20	51.51#
42.00	18.00	37.13#

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Instrument :
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Manual IntegrationsAPPROVED

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.928	152	114594	20.000	ng/ul	0.00
20) Naphthalene-d8	10.728	136	474755	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.557	164	317343	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.316	188	682043	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.398	240	664244	20.000	ng/ul	# 0.00
88) Perylene-d12	23.845	264	648477	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.358	96	14171m	5.393	ng/uL	0.00
4) Pyridine-d5	3.781	84	66409	8.983	ng/ul	0.00
7) Phenol-d5	7.087	99	58333	5.596	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.252	67	188483	30.136	ng/ul	0.00
11) 2-Chlorophenol-d4	7.458	132	173878	22.525	ng/ul	0.00
15) 4-Methylphenol-d8	8.634	113	114794	13.292	ng/ul	0.00
21) Nitrobenzene-d5	9.081	128	113042	29.937	ng/ul	0.00
24) 2-Nitrophenol-d4	9.810	143	119719	29.031	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.352	165	197115	24.553	ng/ul	0.00
31) 4-Chloroaniline-d4	10.863	131	271539	23.589	ng/ul	0.00
46) Dimethylphthalate-d6	13.969	166	765687	31.017	ng/ul	0.00
49) Acenaphthylene-d8	14.251	160	849859	28.446	ng/ul	0.00
54) 4-Nitrophenol-d4	14.763	143	21171m	4.659	ng/ul	0.01
60) Fluorene-d10	15.551	176	665504	31.340	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.675	200	113634	26.349	ng/ul	0.00
73) Anthracene-d10	17.416	188	1137964	34.487	ng/ul	0.00
81) Pyrene-d10	19.645	212	1367562	36.202	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.686	264	1210591	35.609	ng/ul	0.00

Target Compounds Qvalue

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

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