

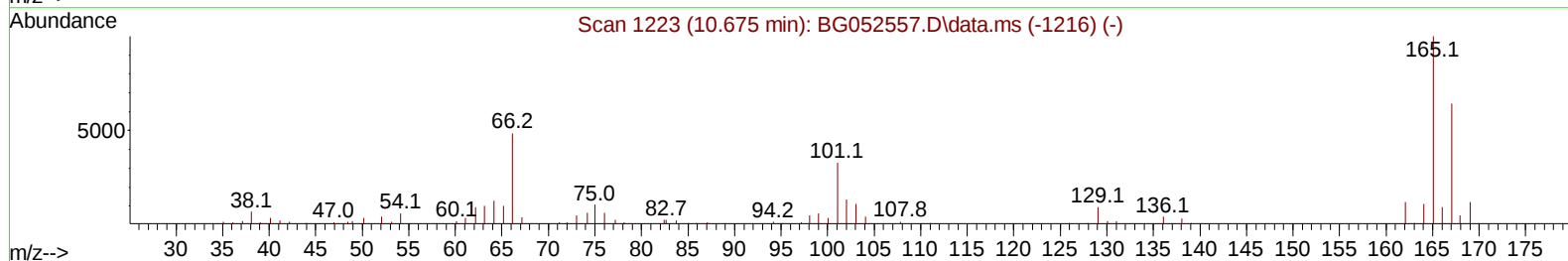
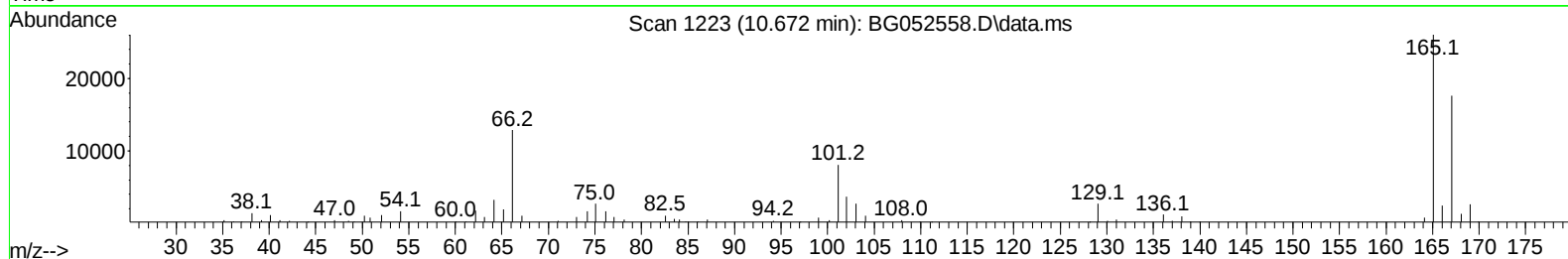
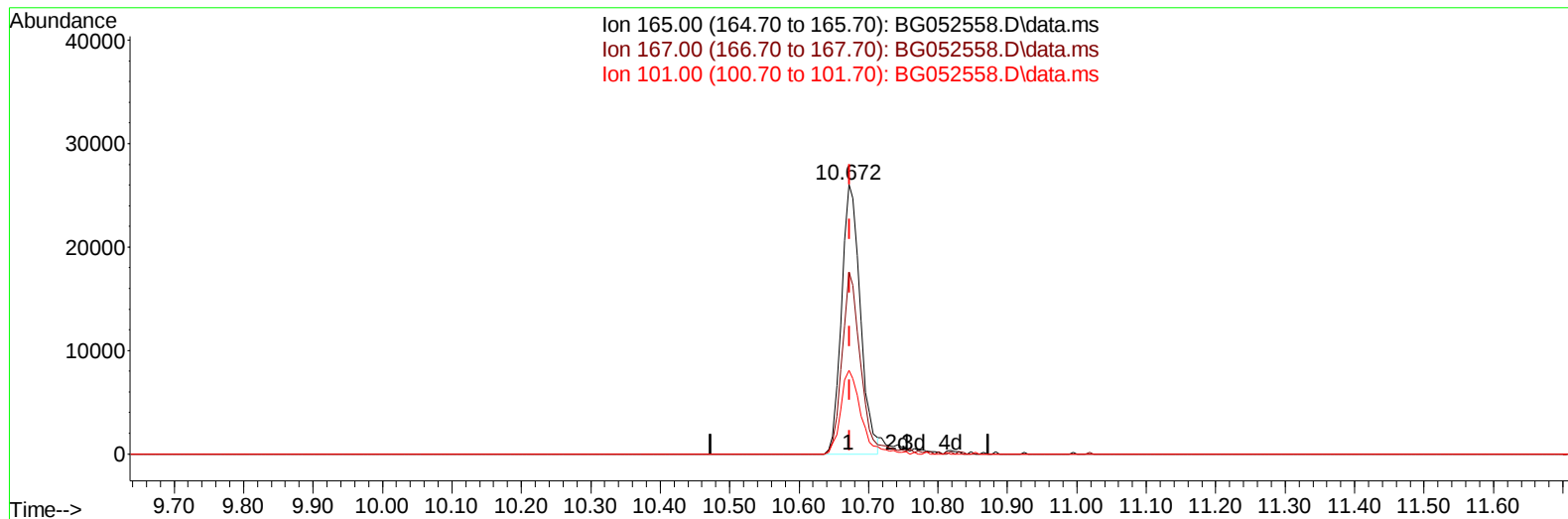
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG030322\
Data File : BG052558.D
Acq On : 3 Mar 2022 17:23
Operator : CG/JU
Sample : PB143054BL
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
SBLK054

Manual IntegrationsAPPROVED

Quant Time: Mar 04 00:02:17 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG022222.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue Feb 22 17:32:12 2022
Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 03/04/2022
Supervised By :Yogesh Patel 03/07/2022



TIC: BG052558.D\data.ms

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

10.672min (-0.001) 24.34 ng/u1

response 48689

Ion	Exp%	Act%
165.00	100.00	100.00
167.00	64.30	67.57
101.00	34.40	31.00
0.00	0.00	0.00

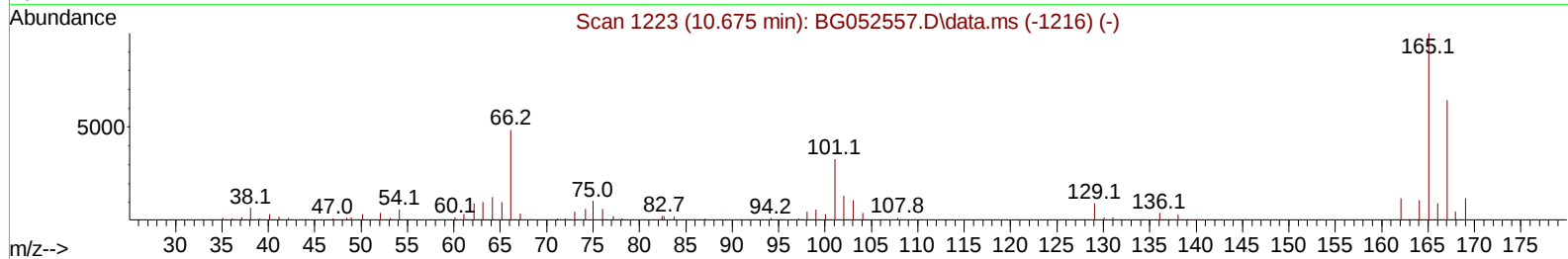
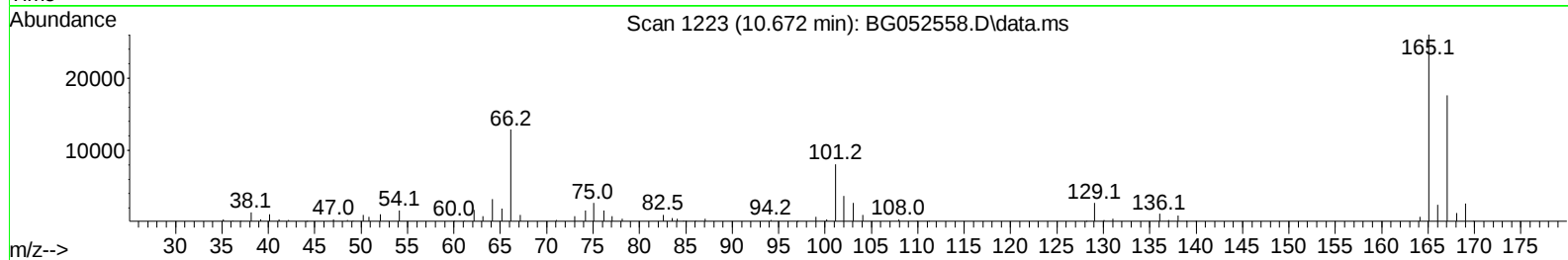
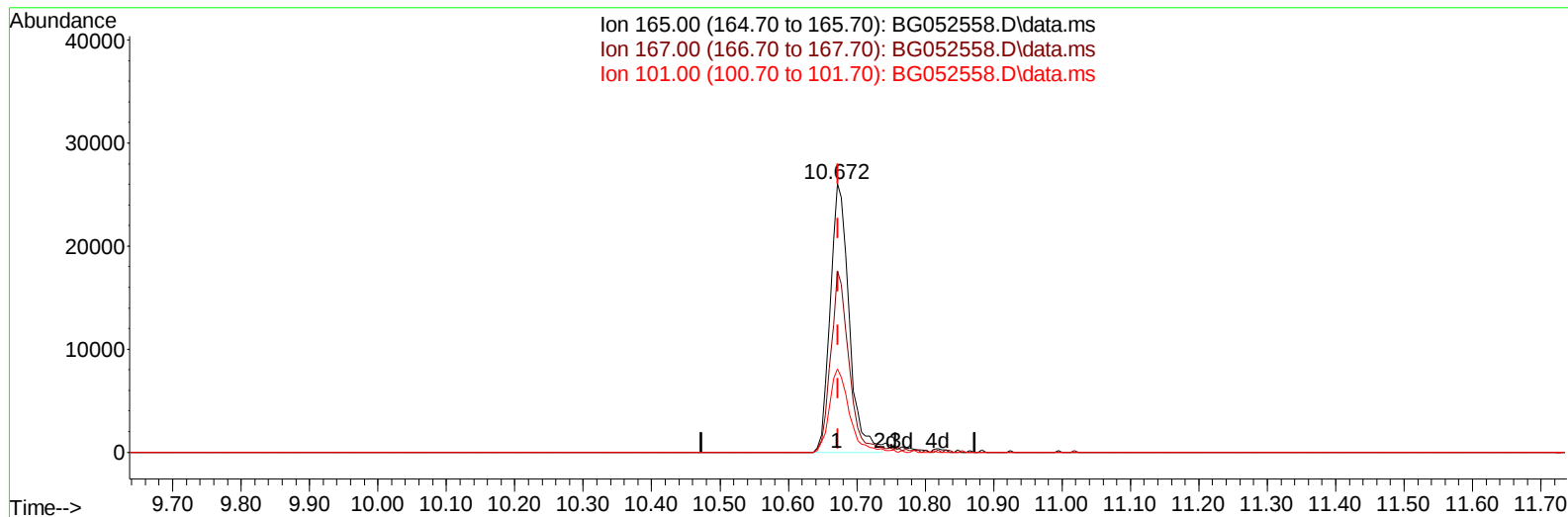
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TIC: BG052558.D\data.ms

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

10.672min (-0.001) 25.03 ng/ul m

response 50086

Ion	Exp%	Act%
165.00	100.00	100.00
167.00	64.30	67.57
101.00	34.40	31.00
0.00	0.00	0.00

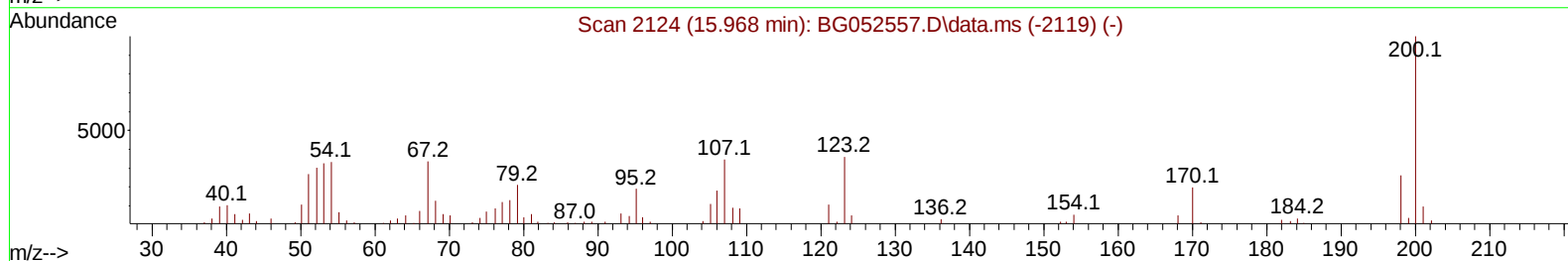
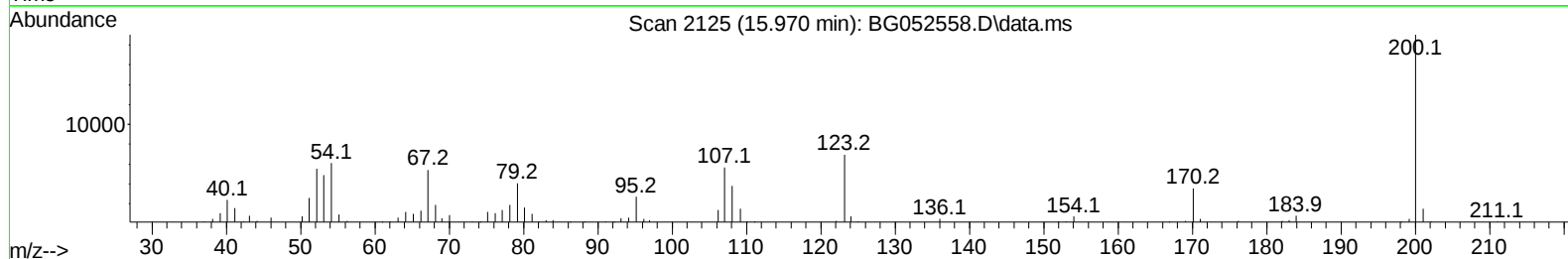
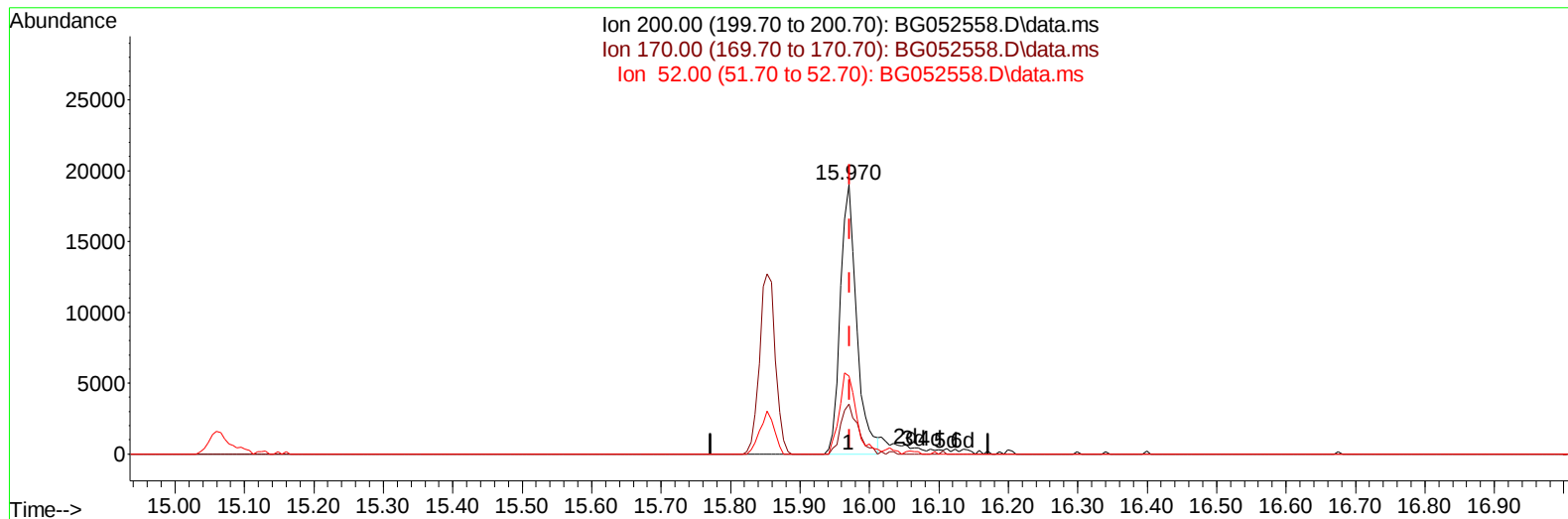
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Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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TIC: BG052558.D\data.ms

(65) 4,6-Dinitro-2-methylphenol-d2 (S)

15.970min (-0.001) 21.67 ng/u1

response 31015

Ion	Exp%	Act%
200.00	100.00	100.00
170.00	19.80	18.49
52.00	47.40	28.97#
0.00	0.00	0.00

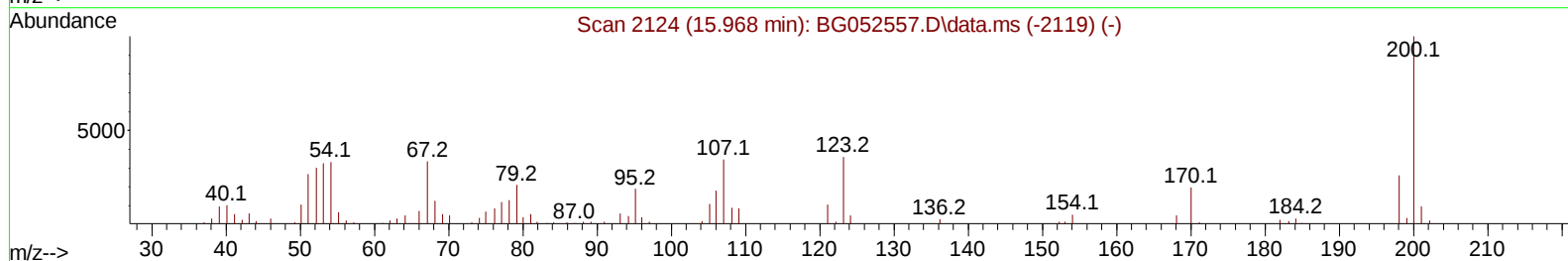
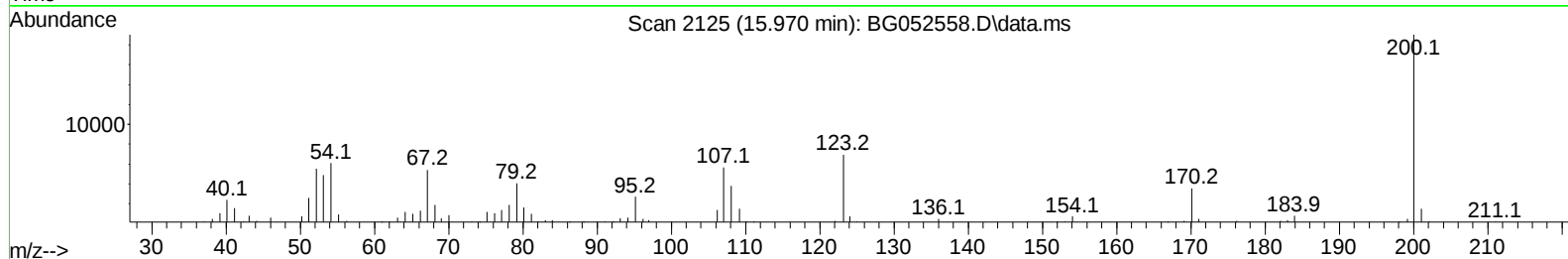
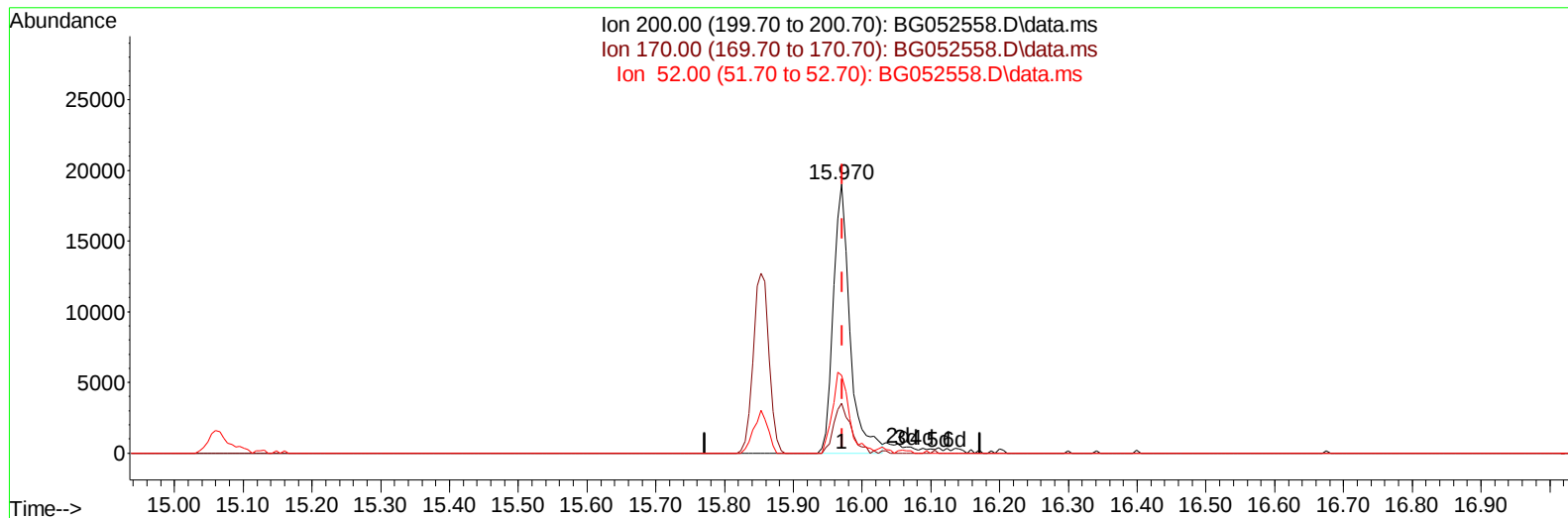
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Acq On : 3 Mar 2022 17:23
Operator : CG/JU
Sample : PB143054BL
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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Manual IntegrationsAPPROVED

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TIC: BG052558.D\data.ms

(65) 4,6-Dinitro-2-methylphenol-d2 (S)

15.970min (-0.001) 22.33 ng/ul m

response 31964

Ion	Exp%	Act%
200.00	100.00	100.00
170.00	19.80	18.49
52.00	47.40	28.97#
0.00	0.00	0.00

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.216	152	26034	20.000	ng/ul	0.00
20) Naphthalene-d8	11.053	136	114898	20.000	ng/ul	# 0.00
38) Acenaphthene-d10	14.860	164	88405	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.615	188	225720	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.927	240	232525	20.000	ng/ul	# 0.00
88) Perylene-d12	25.393	264	231230	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.540	96	3799	5.724	ng/uL	0.00
4) Pyridine-d5	3.975	84	46834	23.392	ng/ul	0.00
7) Phenol-d5	7.364	99	65645	26.592	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.523	67	40750	28.246	ng/ul	0.00
11) 2-Chlorophenol-d4	7.746	132	47076	27.150	ng/ul	0.00
15) 4-Methylphenol-d8	8.927	113	53525	27.867	ng/ul	0.00
21) Nitrobenzene-d5	9.391	128	27010	27.575	ng/ul	0.00
24) 2-Nitrophenol-d4	10.125	143	29739	27.719	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.672	165	50086m	25.034	ng/ul	0.00
31) 4-Chloroaniline-d4	11.183	131	73306	26.439	ng/ul	0.00
46) Dimethylphthalate-d6	14.255	166	214868	29.898	ng/ul	0.00
49) Acenaphthylene-d8	14.560	160	254193	28.652	ng/ul	0.00
54) 4-Nitrophenol-d4	15.060	143	28531	24.973	ng/ul	0.00
60) Fluorene-d10	15.853	176	181873	28.807	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.970	200	31964m	22.333	ng/ul	0.00
73) Anthracene-d10	17.709	188	327905	30.070	ng/ul	0.00
81) Pyrene-d10	19.994	212	424180	31.176	ng/ul	0.00
92) Benzo(a)pyrene-d12	25.152	264	380716	30.690	ng/ul	0.00

Target Compounds Qvalue

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

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