

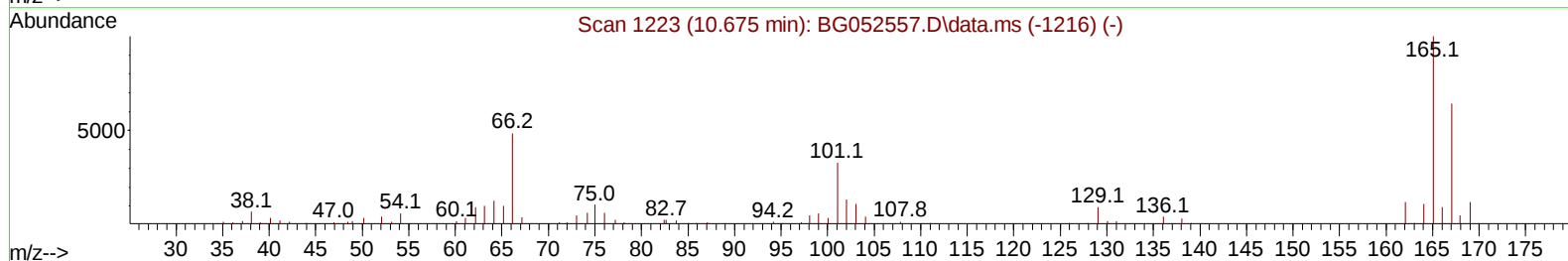
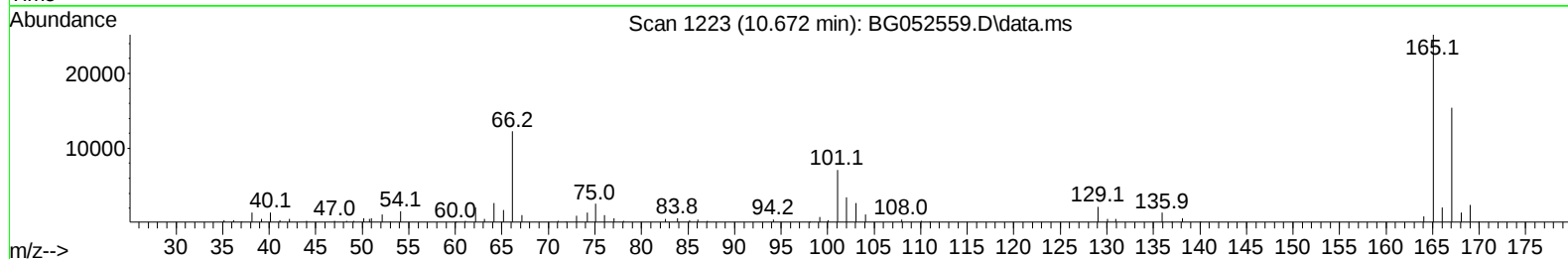
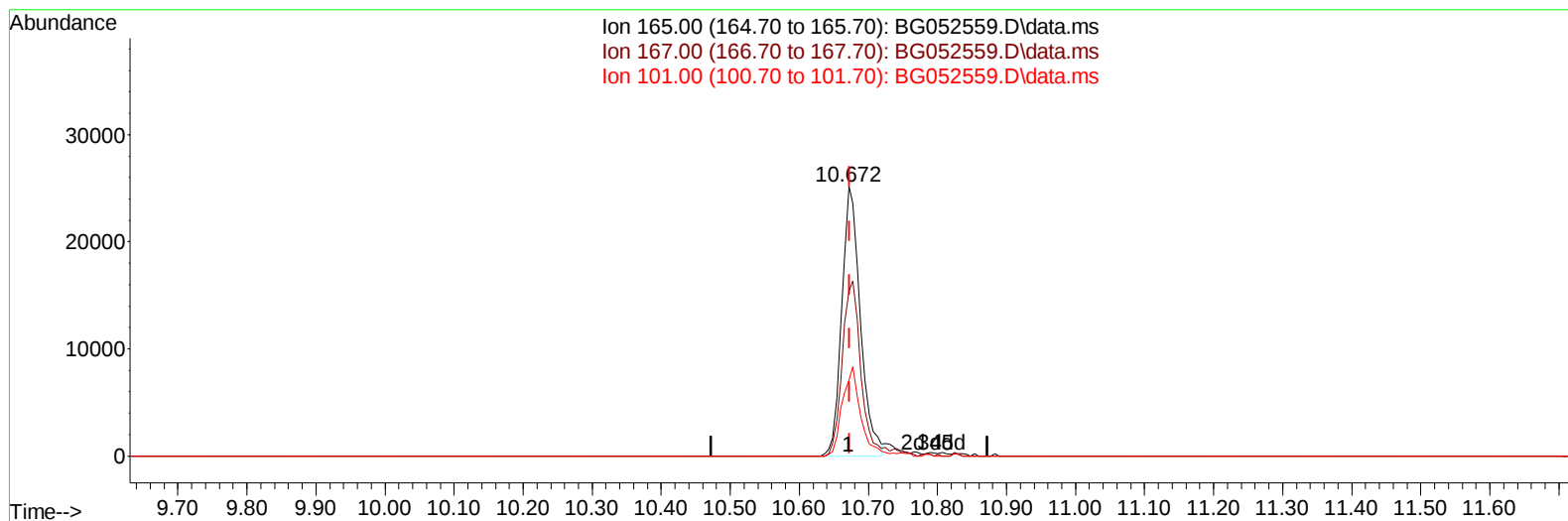
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG030322\
Data File : BG052559.D
Acq On : 3 Mar 2022 18:01
Operator : CG/JU
Sample : PB143070BL
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
SBLK070

Manual IntegrationsAPPROVED

Quant Time: Mar 04 00:02:29 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: BG052559.D\data.ms

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

10.672min (-0.001) 25.34 ng/u1

response 46861

Ion	Exp%	Act%
165.00	100.00	100.00
167.00	64.30	61.30
101.00	34.40	28.29
0.00	0.00	0.00

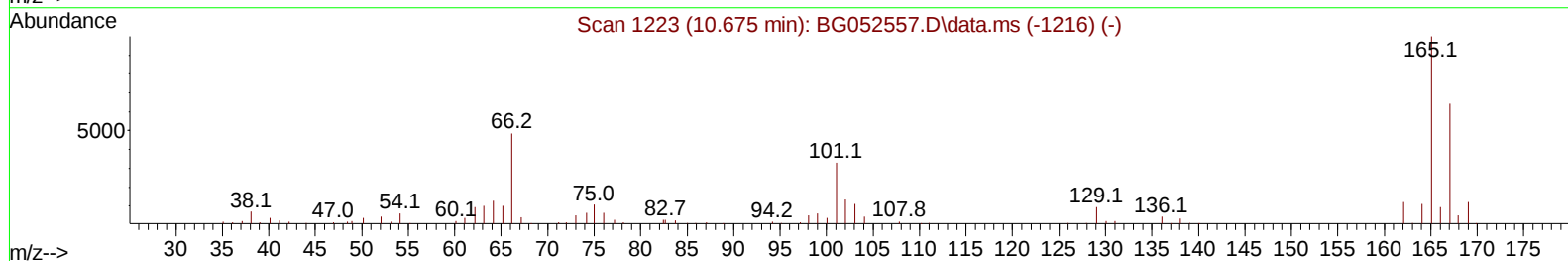
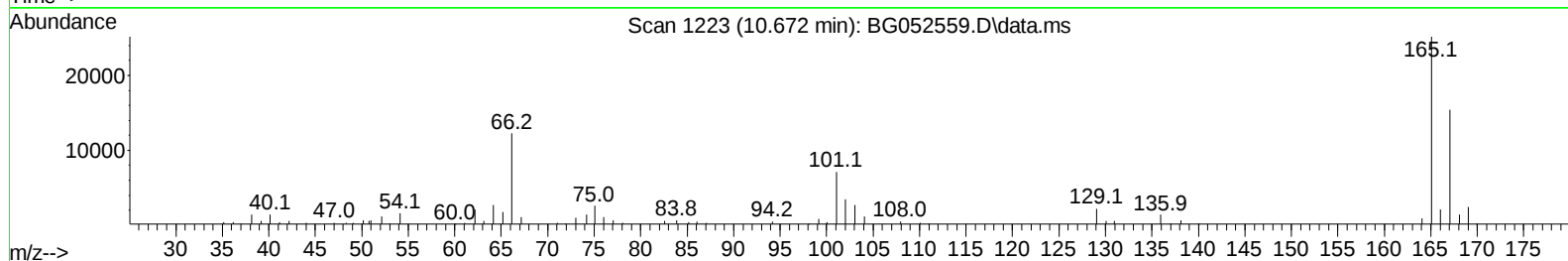
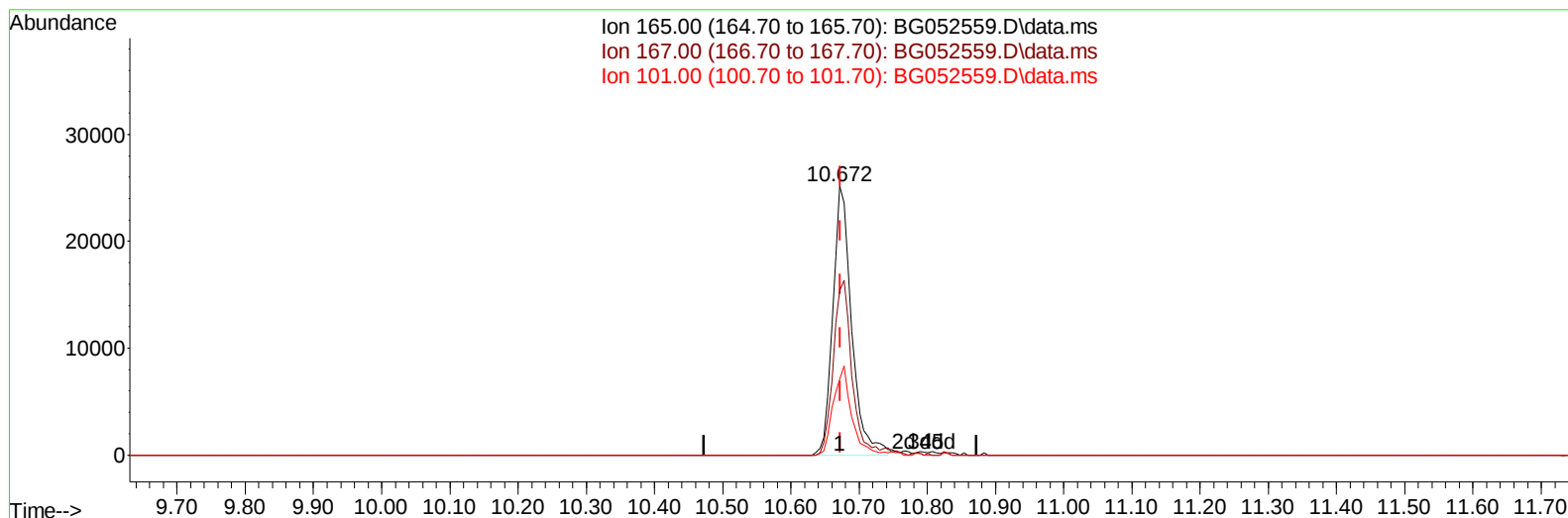
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TIC: BG052559.D\data.ms

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

10.672min (-0.001) 26.03 ng/ul m

response 48139

Ion	Exp%	Act%
165.00	100.00	100.00
167.00	64.30	61.30
101.00	34.40	28.29
0.00	0.00	0.00

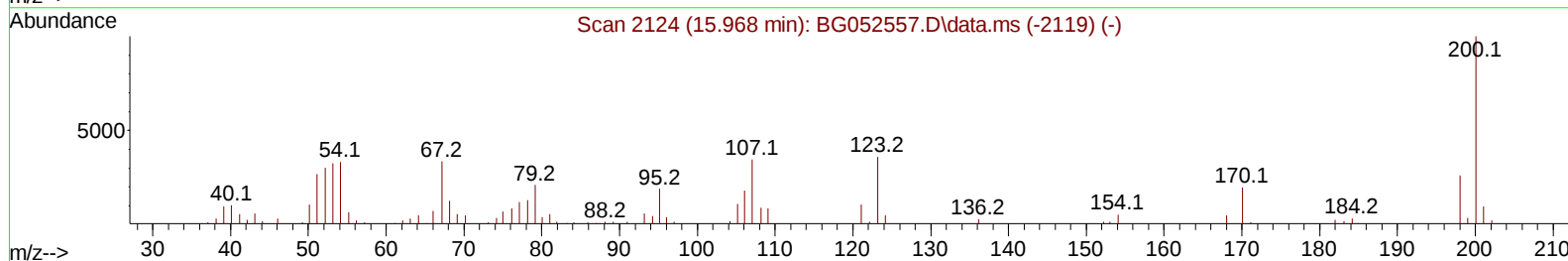
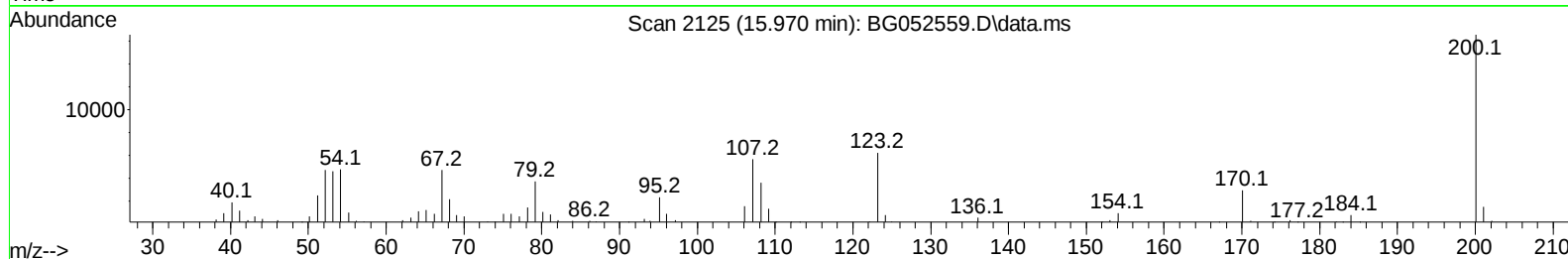
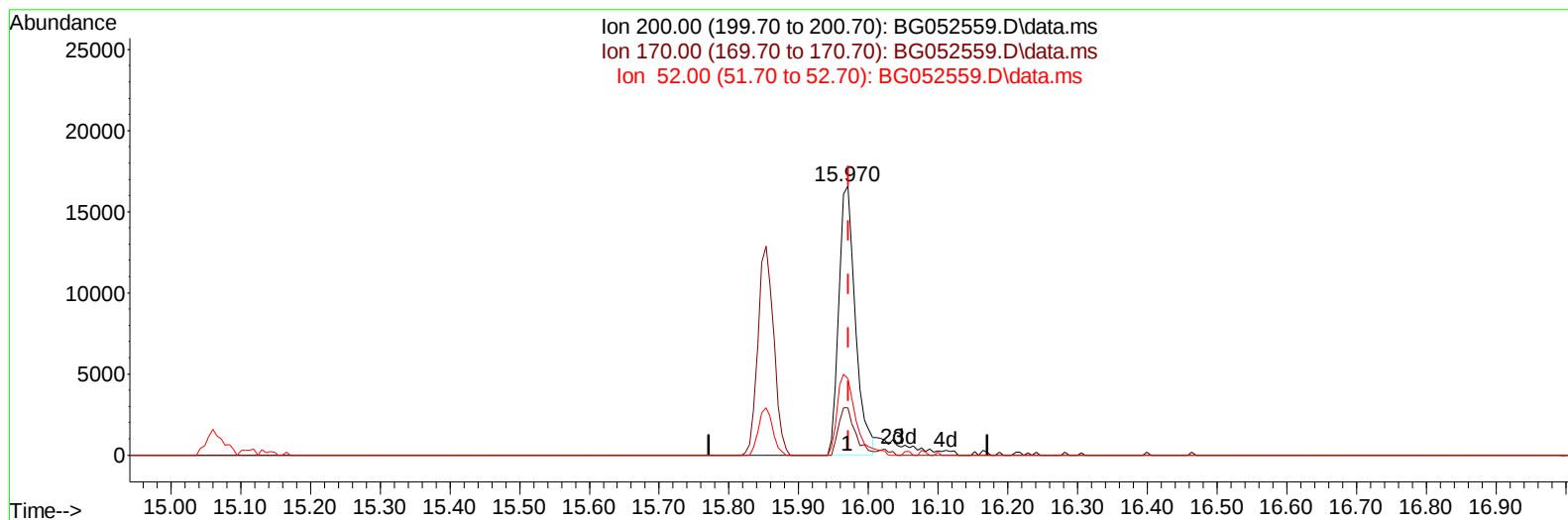
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Manual IntegrationsAPPROVED

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TIC: BG052559.D\data.ms

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

15.970min (-0.001) 20.05 ng/u1

response 27267

Ion	Exp%	Act%
200.00	100.00	100.00
170.00	19.80	17.57
52.00	47.40	28.35#
0.00	0.00	0.00

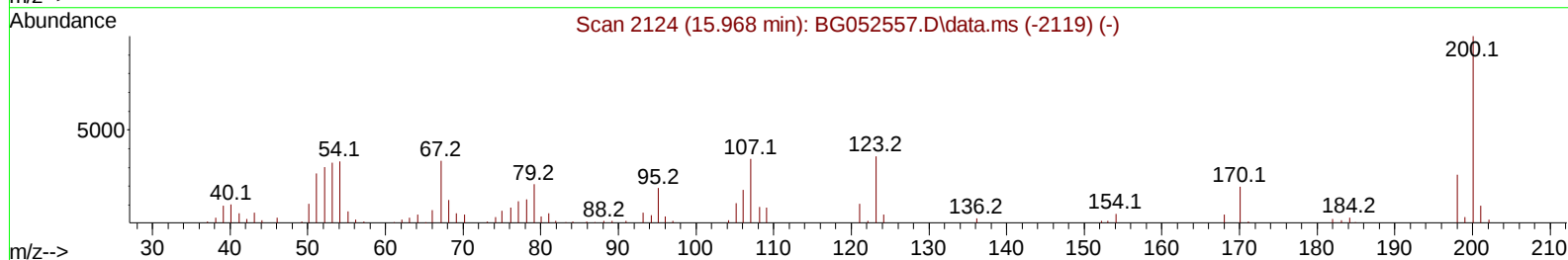
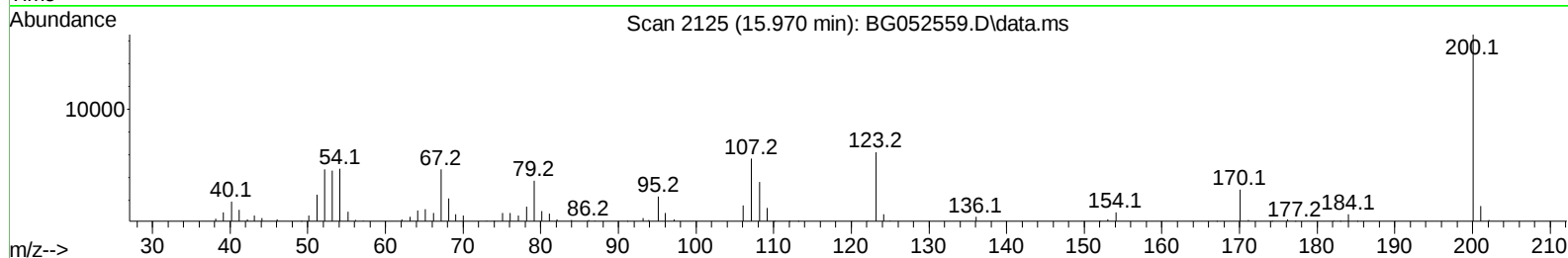
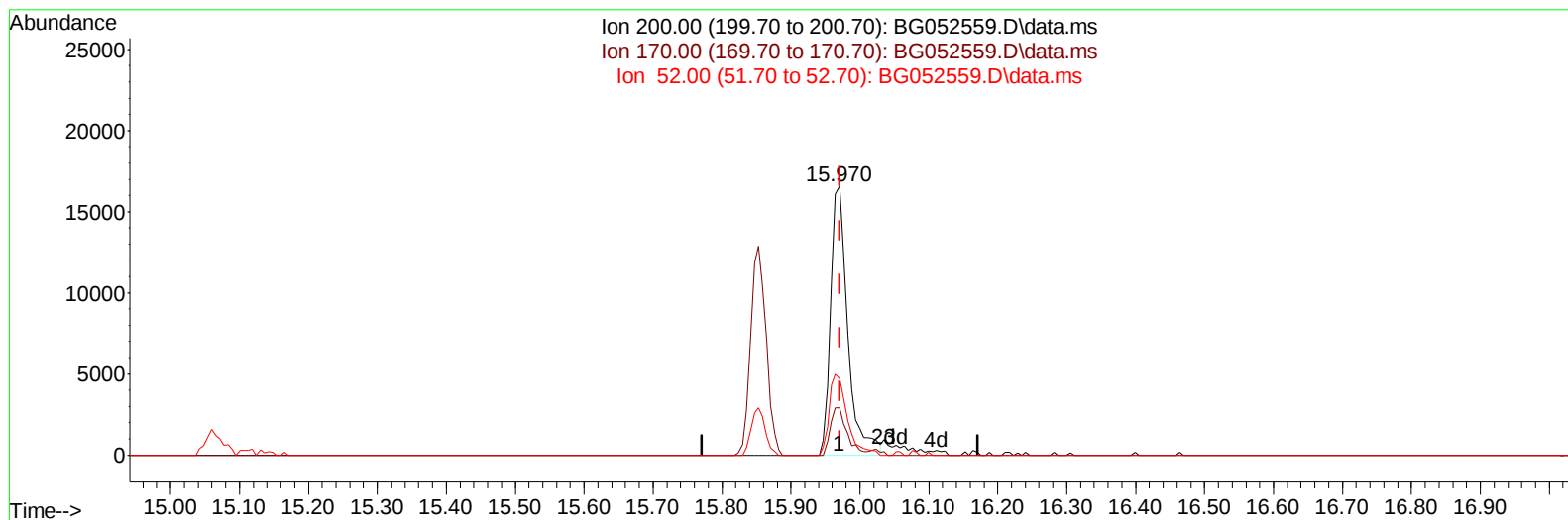
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Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
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TIC: BG052559.D\data.ms

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

15.970min (-0.001) 21.00 ng/ul m

response 28562

Ion	Exp%	Act%
200.00	100.00	100.00
170.00	19.80	17.57
52.00	47.40	28.35#
0.00	0.00	0.00

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.216	152	24532	20.000	ng/ul	0.00
20) Naphthalene-d8	11.054	136	106220	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.866	164	81395	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.615	188	214508	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.927	240	214895	20.000	ng/ul	# 0.00
88) Perylene-d12	25.387	264	214019	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.540	96	3532	5.648	ng/uL	0.00
4) Pyridine-d5	3.969	84	44875	23.786	ng/ul	0.00
7) Phenol-d5	7.364	99	61410	26.399	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.523	67	38809	28.548	ng/ul	0.00
11) 2-Chlorophenol-d4	7.746	132	44474	27.220	ng/ul	0.00
15) 4-Methylphenol-d8	8.927	113	48954	27.047	ng/ul	0.00
21) Nitrobenzene-d5	9.391	128	24630	27.200	ng/ul	0.00
24) 2-Nitrophenol-d4	10.120	143	28217	28.449	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.672	165	48139m	26.026	ng/ul	0.00
31) 4-Chloroaniline-d4	11.183	131	68840	26.857	ng/ul	0.00
46) Dimethylphthalate-d6	14.255	166	205174	31.008	ng/ul	0.00
49) Acenaphthylene-d8	14.561	160	240073	29.390	ng/ul	0.00
54) 4-Nitrophenol-d4	15.060	143	23898	22.720	ng/ul	0.00
60) Fluorene-d10	15.853	176	174443	30.010	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.970	200	28562m	21.000	ng/ul	0.00
73) Anthracene-d10	17.709	188	309654	29.881	ng/ul	0.00
81) Pyrene-d10	19.994	212	394383	31.364	ng/ul	0.00
92) Benzo(a)pyrene-d12	25.146	264	358419	31.216	ng/ul	-0.01

Target Compounds Qvalue

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

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