

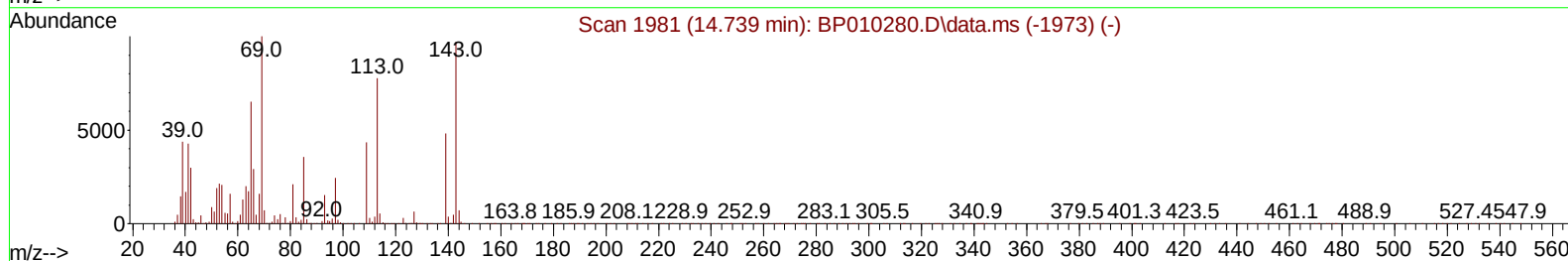
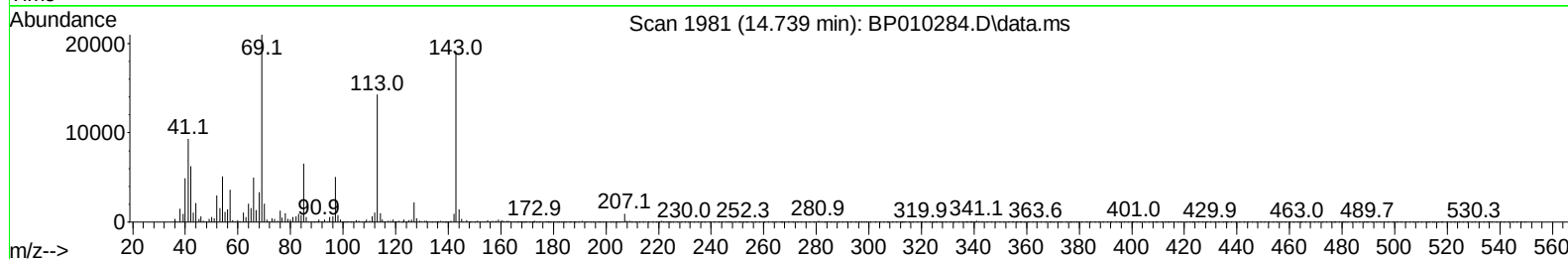
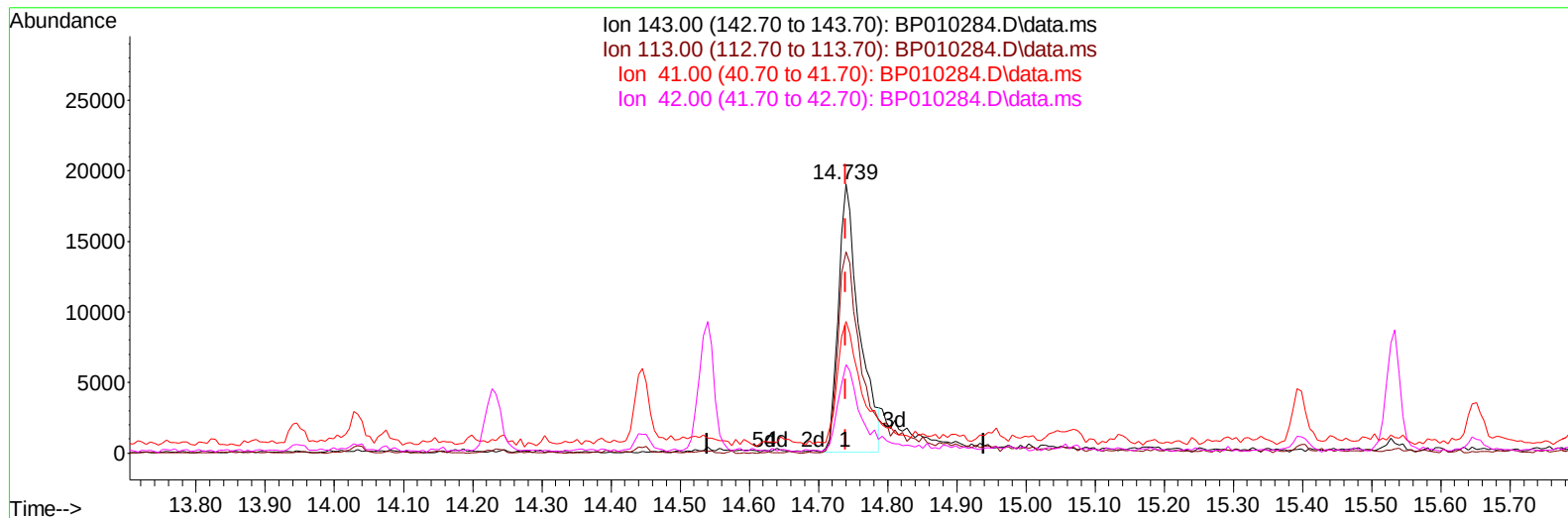
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051122\
Data File : BP010284.D
Acq On : 11 May 2022 12:02
Operator : CG/JU
Sample : N2632-11
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EW4B3

Manual IntegrationsAPPROVED

Quant Time: May 12 01:12:12 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP051022.M
Quant Title : SVOA CALIBRATION
QLast Update : Thu May 12 01:10:15 2022
Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 05/12/2022
Supervised By :Yogesh Patel 05/17/2022



TIC: BP010284.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.739min (-0.000) 6.64 ng/u1

response 39664

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	132.70	74.88#
41.00	23.20	48.96#
42.00	18.00	32.87#

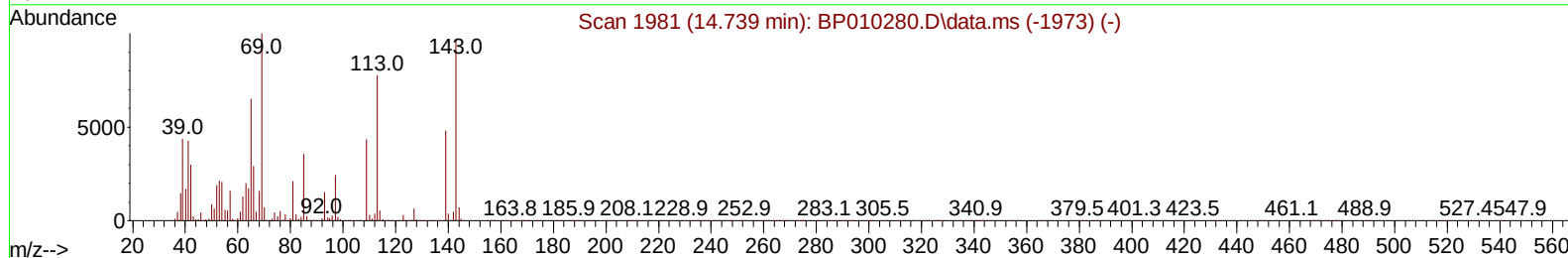
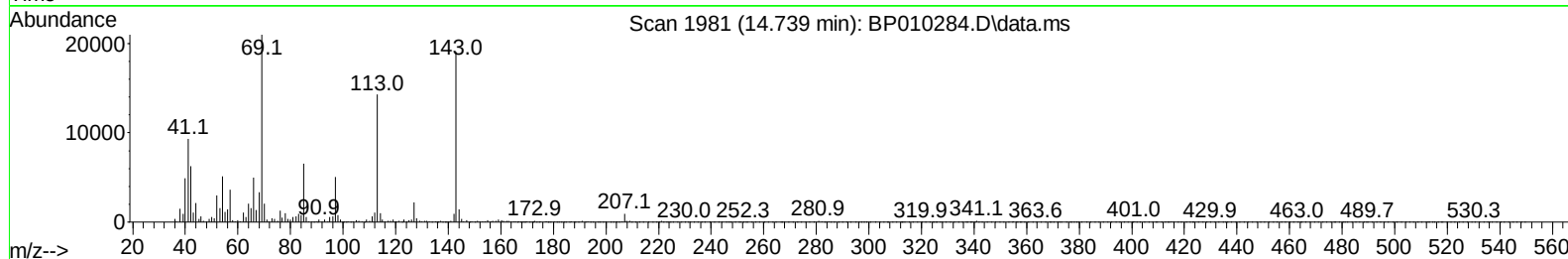
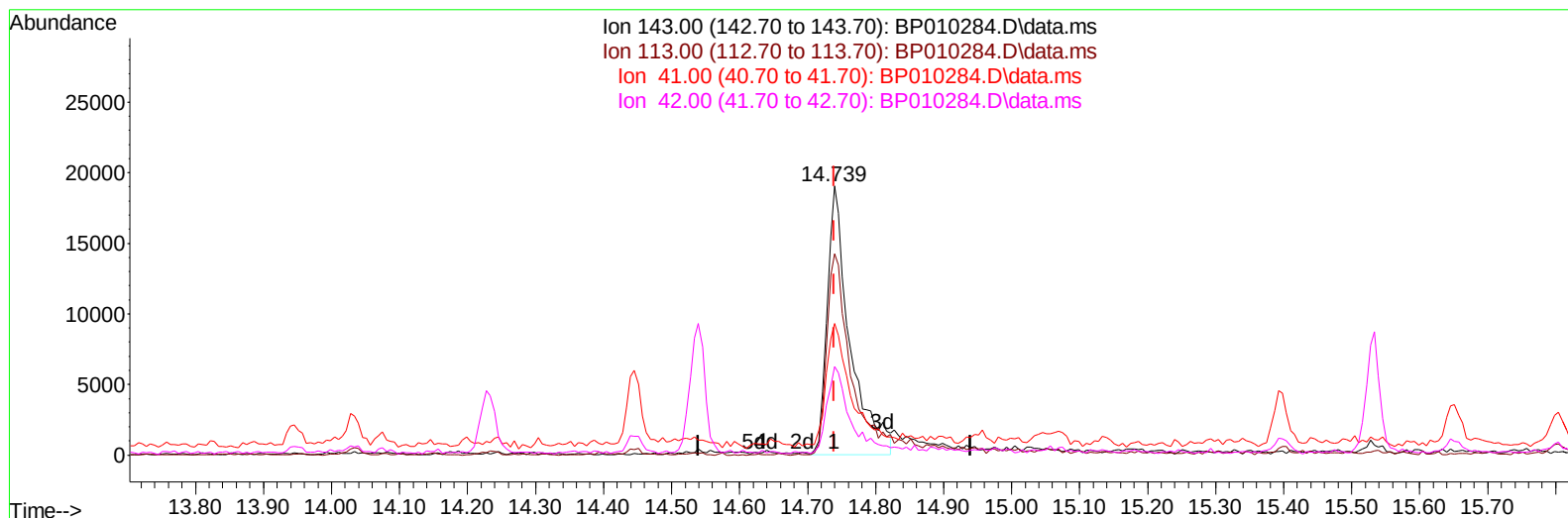
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TIC: BP010284.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.739min (-0.000) 7.44 ng/ul m

response 44397

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	132.70	74.88#
41.00	23.20	48.96#
42.00	18.00	32.87#

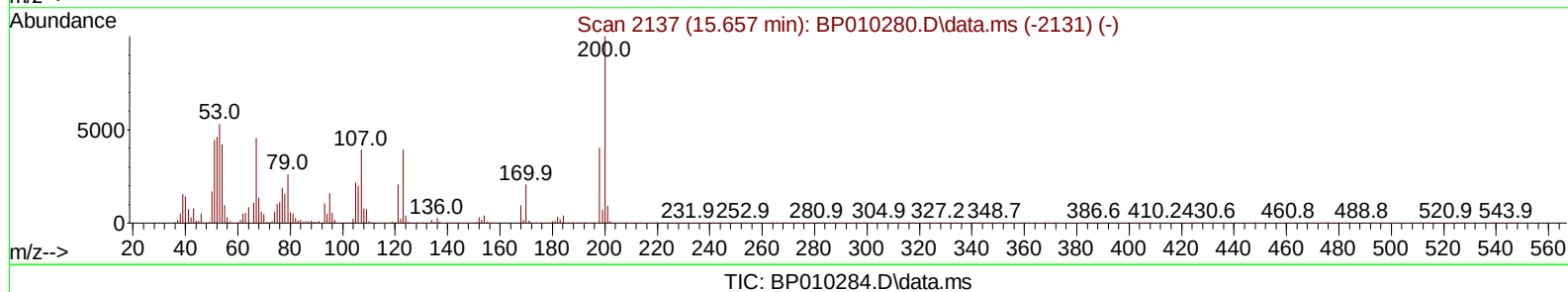
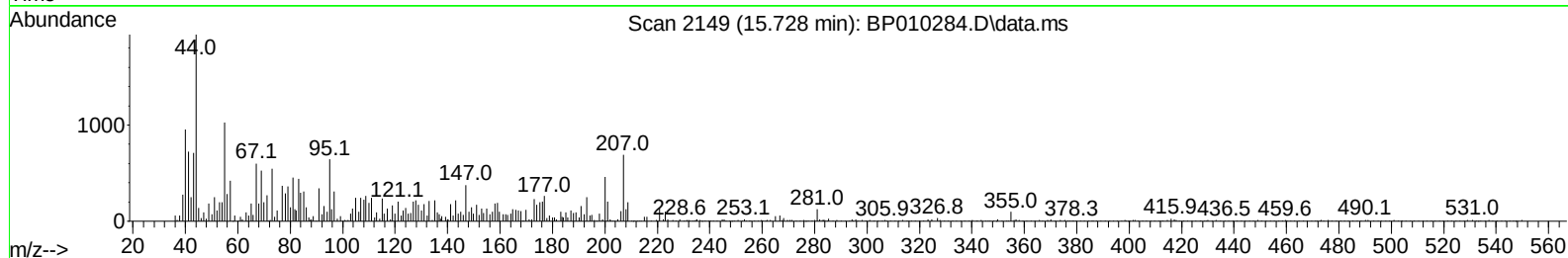
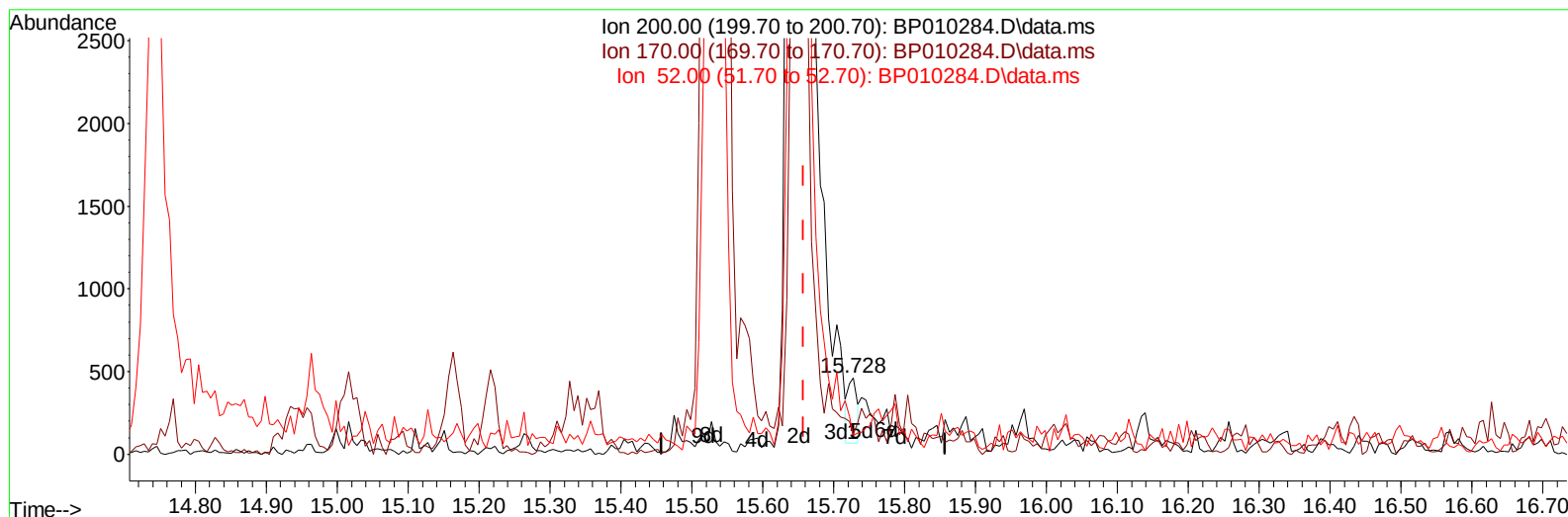
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Instrument :
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(65) 4,6-Dinitro-2-methylphenol-d2 (S)

15.728min (+ 0.071) 0.06 ng/ul

response 350

Ion	Exp%	Act%
200.00	100.00	100.00
170.00	23.00	26.13
52.00	21.20	24.84
0.00	0.00	0.00

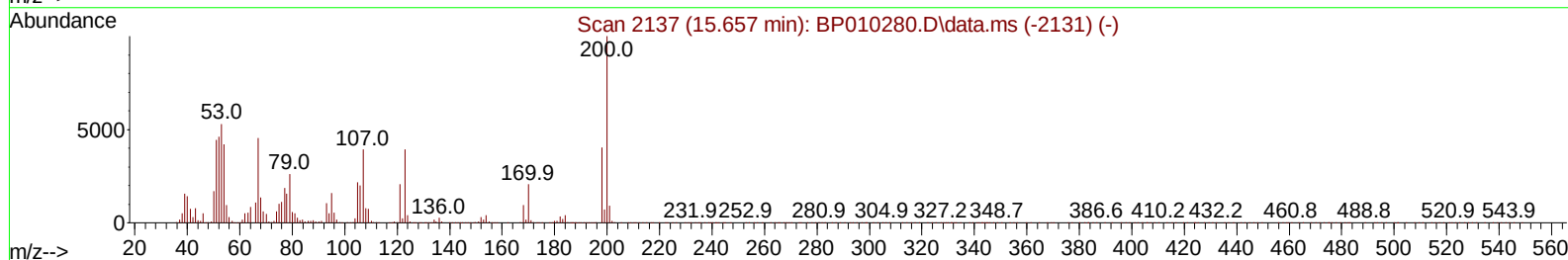
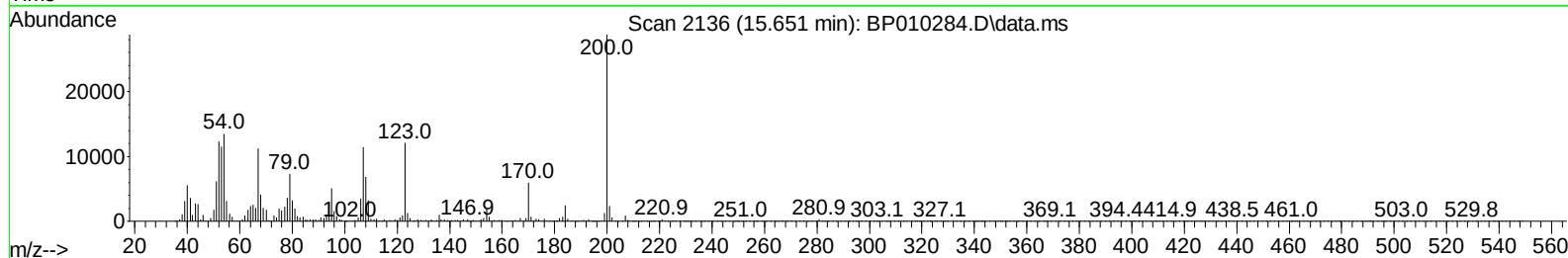
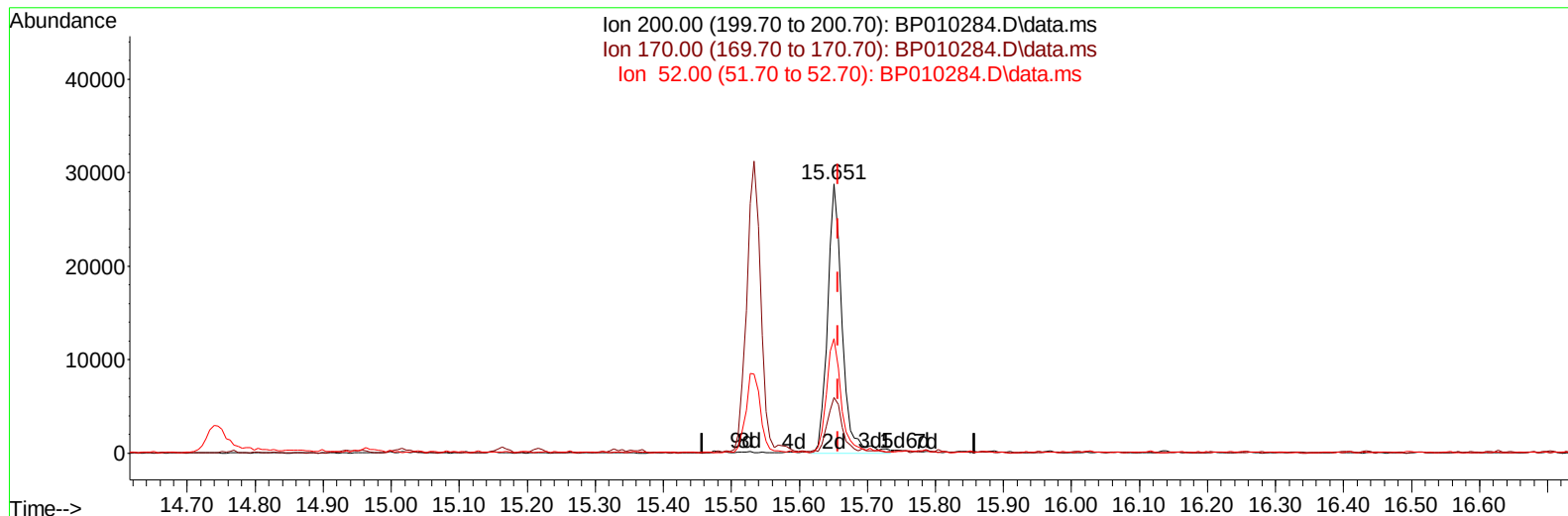
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Instrument :
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TIC: BP010284.D\data.ms

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

15.651min (-0.006) 7.12 ng/ul m

response 42590

Ion	Exp%	Act%
200.00	100.00	100.00
170.00	23.00	20.61
52.00	21.20	42.66#
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	7.910	152	122217	20.000	ng/ul	0.00
20) Naphthalene-d8	10.710	136	545966	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.539	164	390734	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.292	188	873164	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.374	240	907407	20.000	ng/ul	0.00
88) Perylene-d12	23.810	264	956797	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.352	96	7675	2.668	ng/uL	0.00
4) Pyridine-d5	3.787	84	21275	2.657	ng/ul	0.02
7) Phenol-d5	7.069	99	131945	11.624	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.234	67	98485	14.025	ng/ul	0.00
11) 2-Chlorophenol-d4	7.440	132	105240	12.589	ng/ul	0.00
15) 4-Methylphenol-d8	8.616	113	112325	12.036	ng/ul	0.00
21) Nitrobenzene-d5	9.063	128	57718	13.185	ng/ul	0.00
24) 2-Nitrophenol-d4	9.793	143	59480	11.917	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.334	165	106647	11.739	ng/ul	0.00
31) 4-Chloroaniline-d4	10.846	131	154543	11.523	ng/ul	0.00
46) Dimethylphthalate-d6	13.945	166	442228	14.207	ng/ul	0.00
49) Acenaphthylene-d8	14.234	160	473495	12.758	ng/ul	0.00
54) 4-Nitrophenol-d4	14.739	143	44397m	7.436	ng/ul	0.00
60) Fluorene-d10	15.534	176	382966	14.575	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.651	200	42590m	7.123	ng/ul	0.00
73) Anthracene-d10	17.386	188	640258	14.974	ng/ul	0.00
81) Pyrene-d10	19.622	212	870759	17.365	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.651	264	864163	16.936	ng/ul	0.00

Target Compounds Qvalue

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

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