

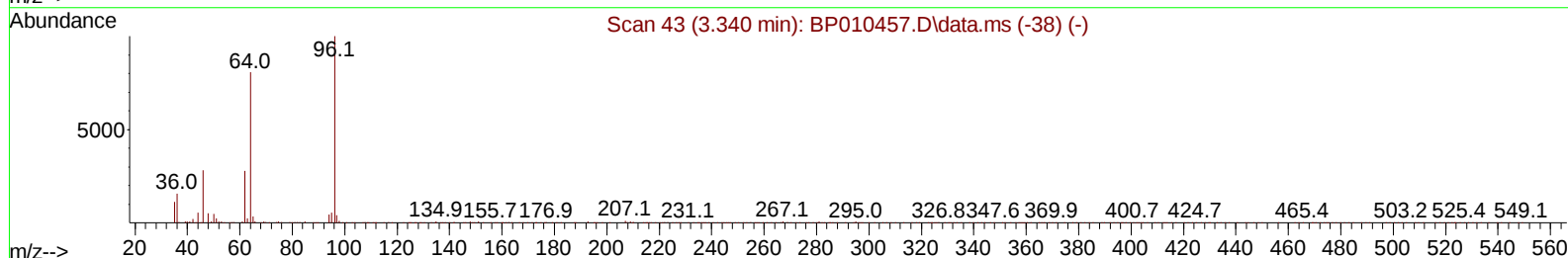
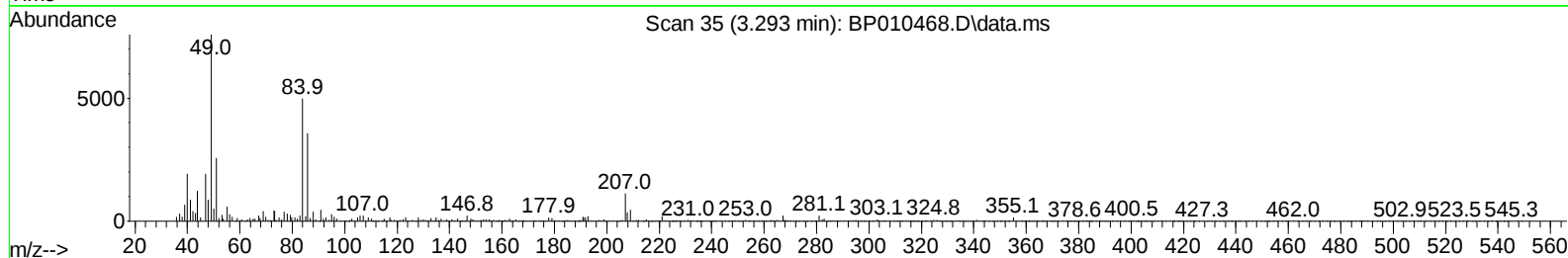
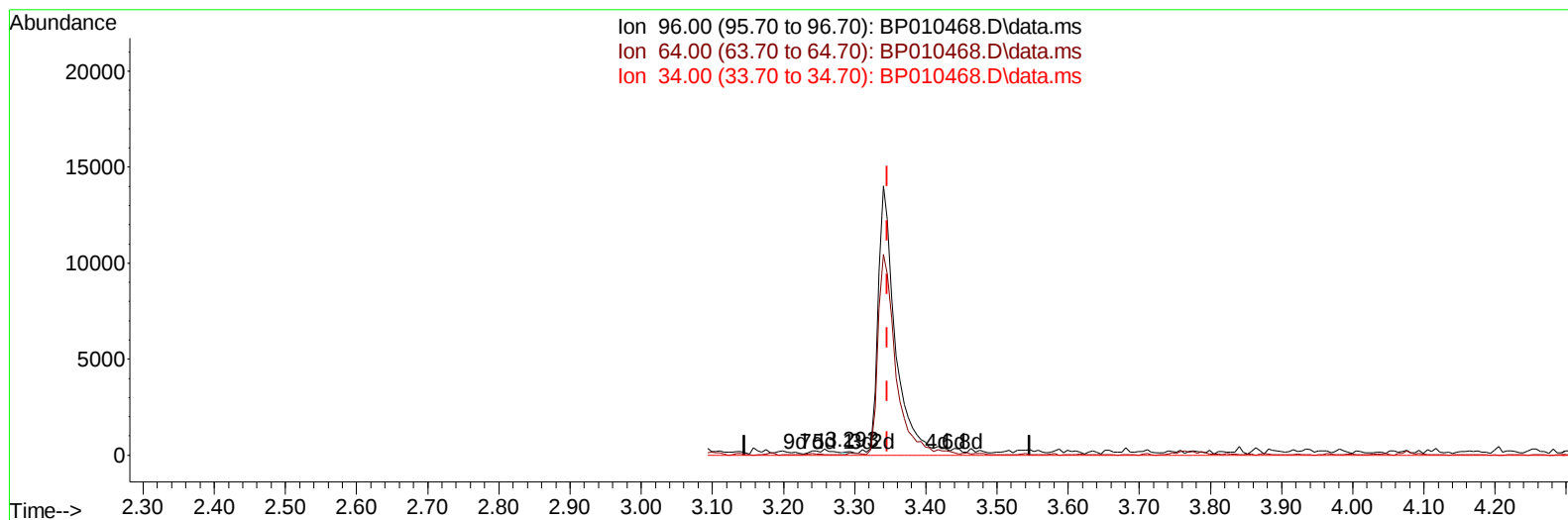
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP052022\
 Data File : BP010468.D
 Acq On : 21 May 2022 10:23
 Operator : CG/JU
 Sample : N2918-23
 Misc :
 ALS Vial : 43 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 DBTD9

Manual IntegrationsAPPROVED

Quant Time: May 23 00:44:07 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP051322.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri May 13 15:04:31 2022
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 05/23/2022
 Supervised By :mohammad ahmed 05/25/2022



TIC: BP010468.D\data.ms

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

3.293min (-0.053) 0.02 ng/uL

response 86

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	64.86
34.00	0.00	0.00
0.00	0.00	0.00

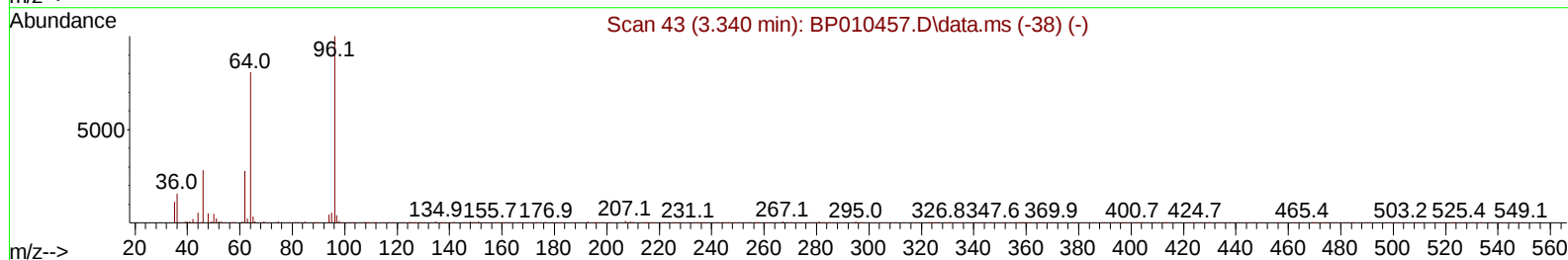
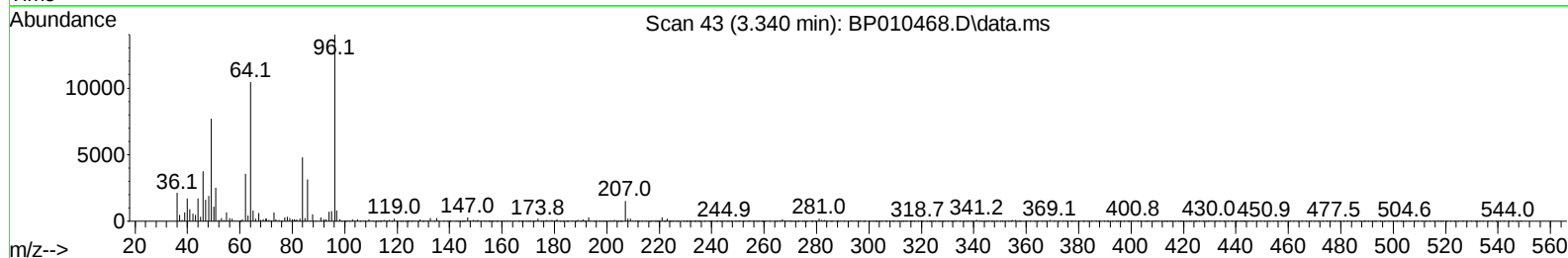
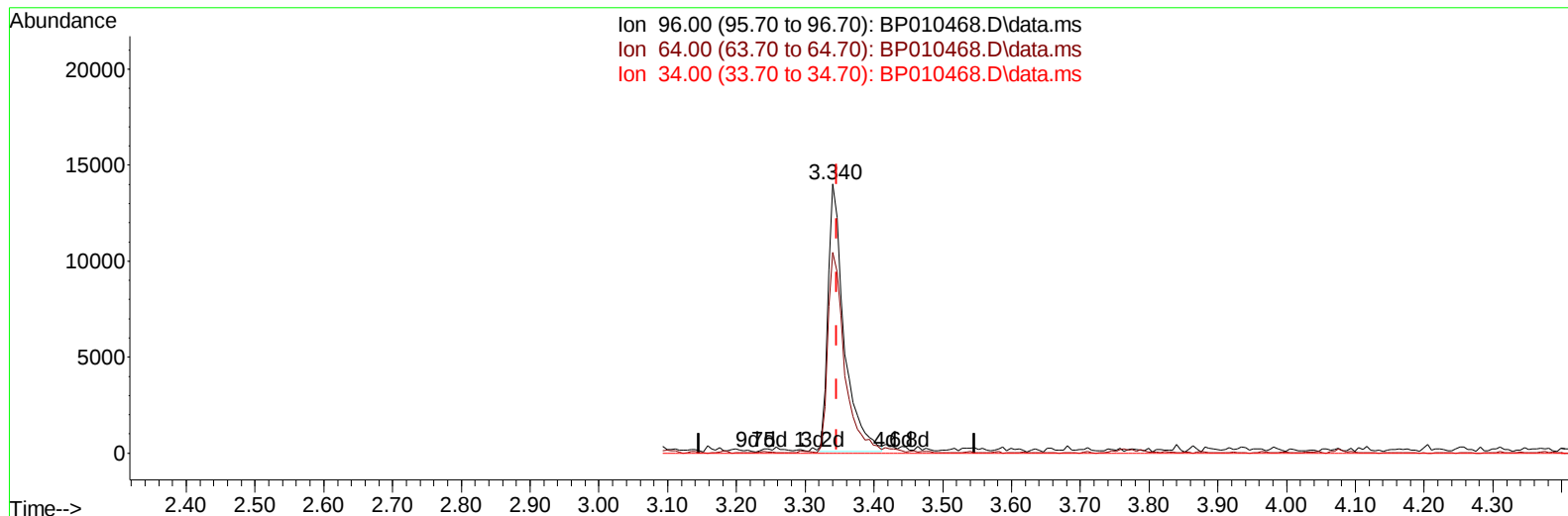
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TIC: BP010468.D\data.ms

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

3.340min (-0.006) 4.65 ng/uL m

response 22812

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	74.67#
34.00	0.00	0.00
0.00	0.00	0.00

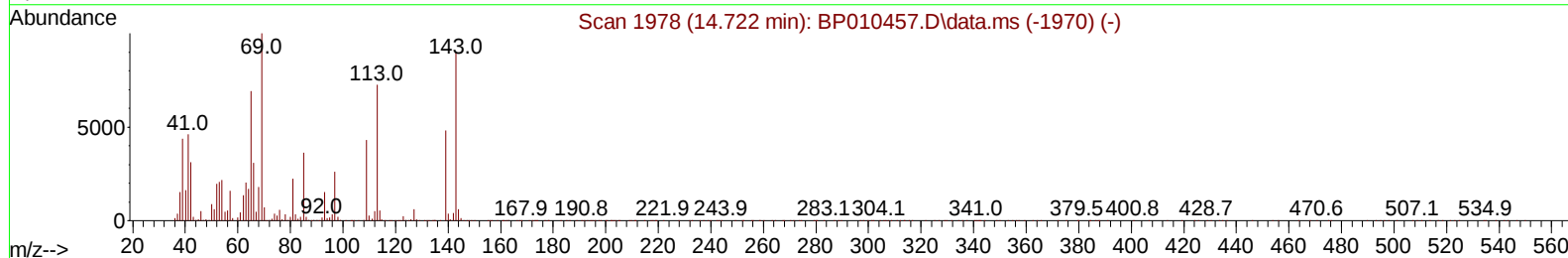
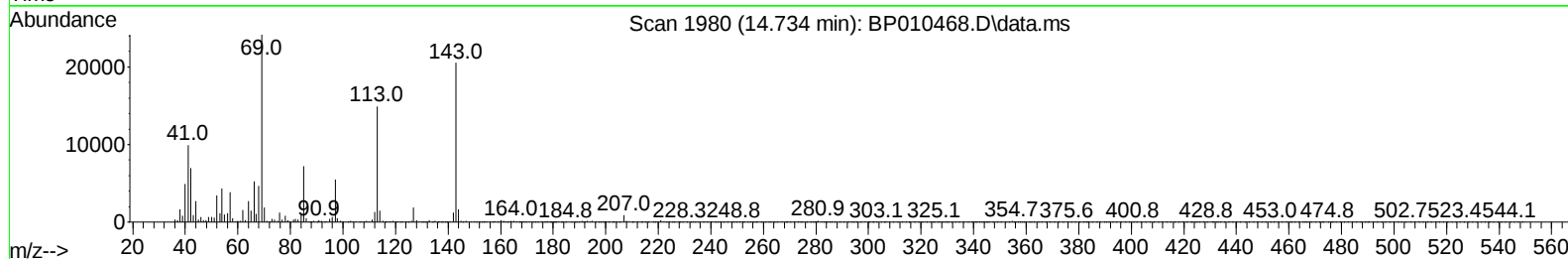
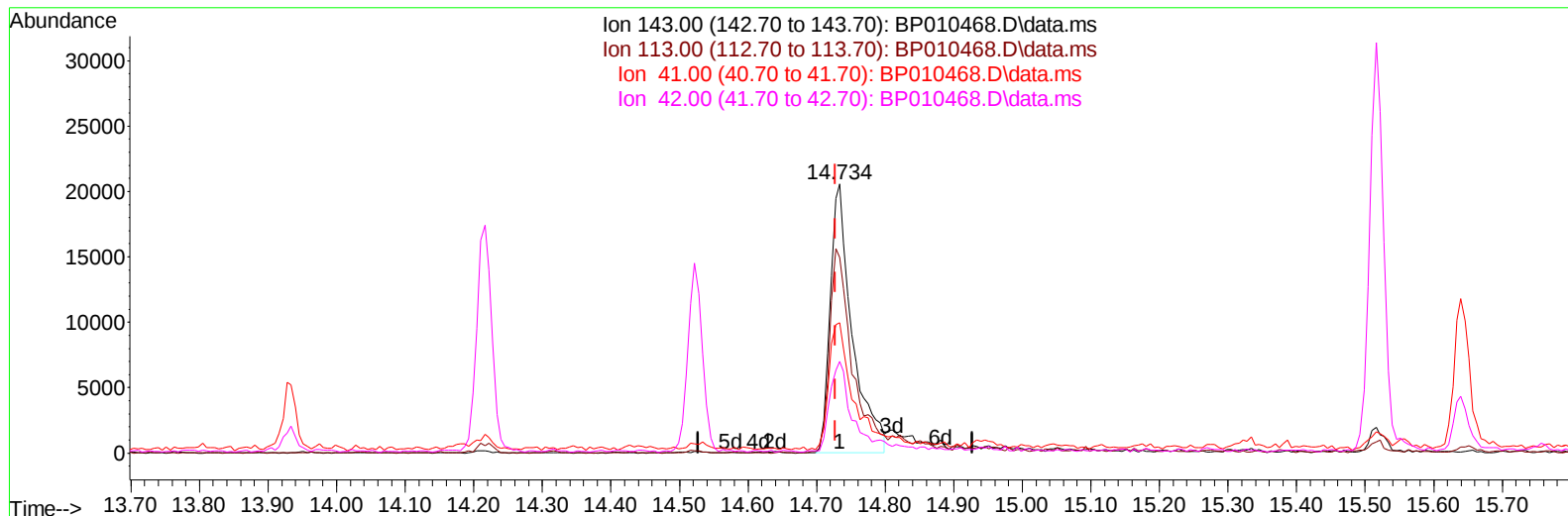
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TIC: BP010468.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.734min (+ 0.006) 5.67 ng/ul

response 45802

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	132.70	72.72#
41.00	23.20	48.39#
42.00	18.00	34.07#

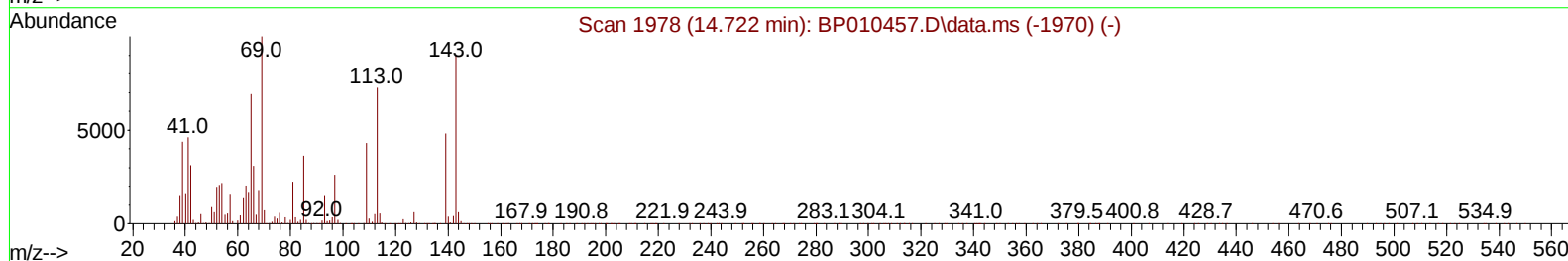
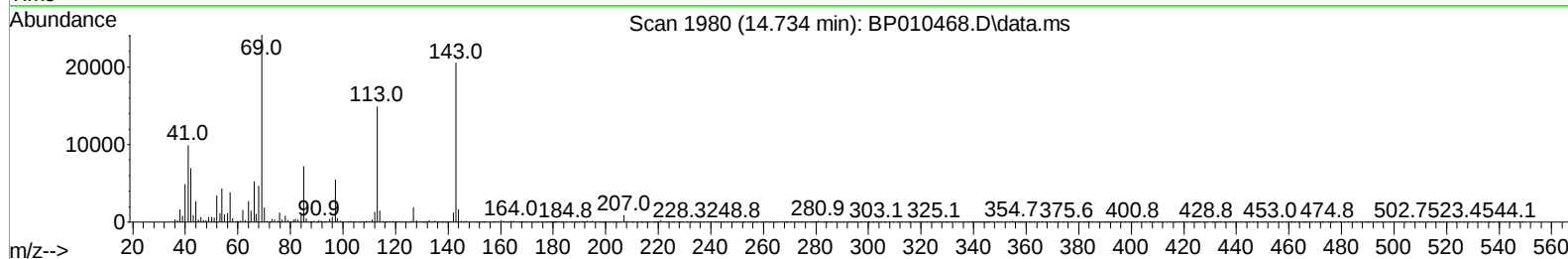
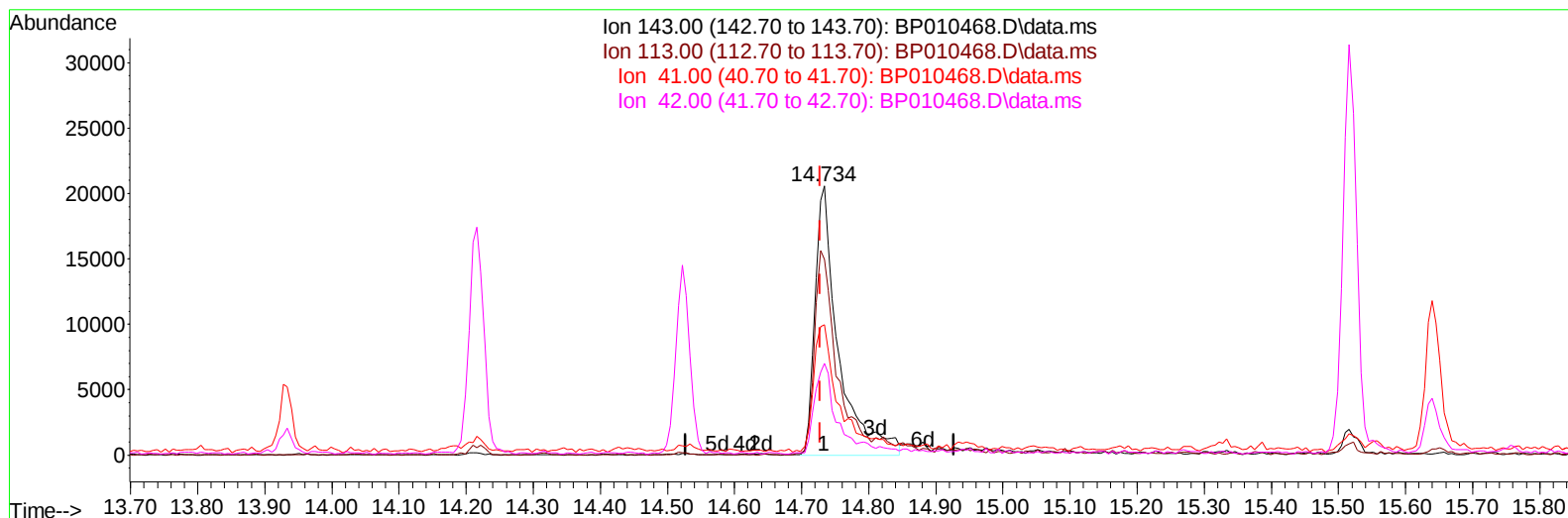
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Instrument :
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TIC: BP010468.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.734min (+ 0.006) 6.16 ng/ul m

response 49774

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	132.70	72.72#
41.00	23.20	48.39#
42.00	18.00	34.07#

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.893	152	209207	20.000	ng/ul	0.00
20) Naphthalene-d8	10.693	136	934752	20.000	ng/ul	-0.01
38) Acenaphthene-d10	14.522	164	615983	20.000	ng/ul	-0.01
64) Phenanthrene-d10	17.275	188	1288509	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.357	240	1037302	20.000	ng/ul	# 0.00
88) Perylene-d12	23.780	264	830918	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.340	96	22812m	4.648	ng/uL	0.00
4) Pyridine-d5	3.764	84	101193	7.261	ng/ul	0.00
7) Phenol-d5	7.058	99	129098	7.352	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.222	67	427869	36.268	ng/ul	0.00
11) 2-Chlorophenol-d4	7.422	132	386985	28.802	ng/ul	-0.01
15) 4-Methylphenol-d8	8.599	113	267979	17.958	ng/ul	-0.01
21) Nitrobenzene-d5	9.052	128	269610	37.268	ng/ul	0.00
24) 2-Nitrophenol-d4	9.775	143	286753	37.402	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.316	165	464468	32.040	ng/ul	0.00
31) 4-Chloroaniline-d4	10.828	131	516145	23.946	ng/ul	-0.01
46) Dimethylphthalate-d6	13.934	166	1772907	39.065	ng/ul	0.00
49) Acenaphthylene-d8	14.216	160	2036919	38.911	ng/ul	0.00
54) 4-Nitrophenol-d4	14.734	143	49774m	6.162	ng/ul	0.00
60) Fluorene-d10	15.516	176	1545693	39.148	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.640	200	182179	24.175	ng/ul	0.00
73) Anthracene-d10	17.375	188	2558096	43.844	ng/ul	0.00
81) Pyrene-d10	19.610	212	2932358	53.231	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.627	264	1895823	45.438	ng/ul	0.00
Target Compounds						
47) Dimethylphthalate	13.981	163	181972	4.042	ng/ul#	Qvalue 93

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

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