

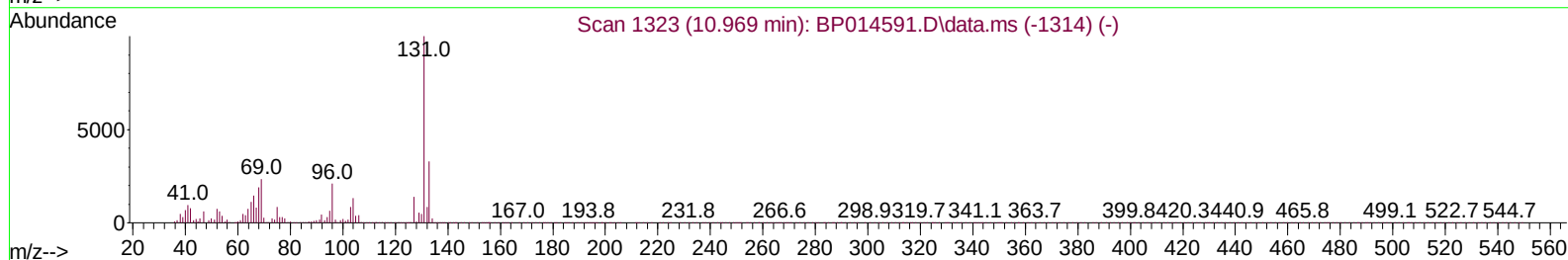
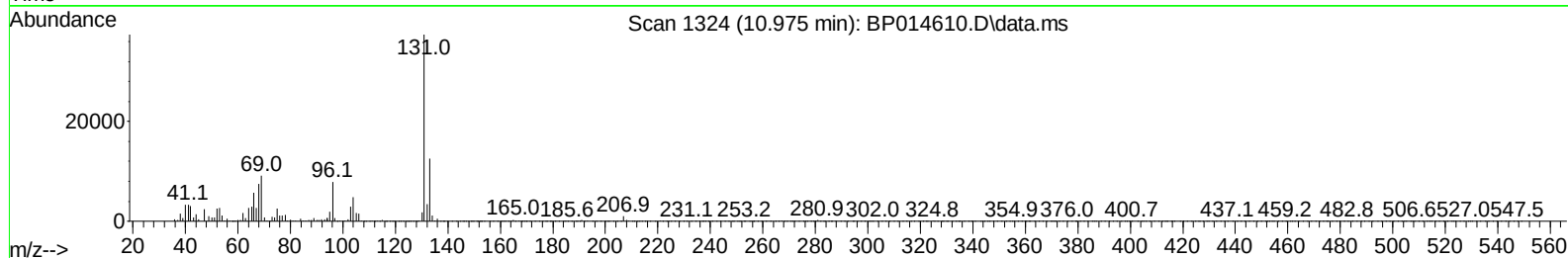
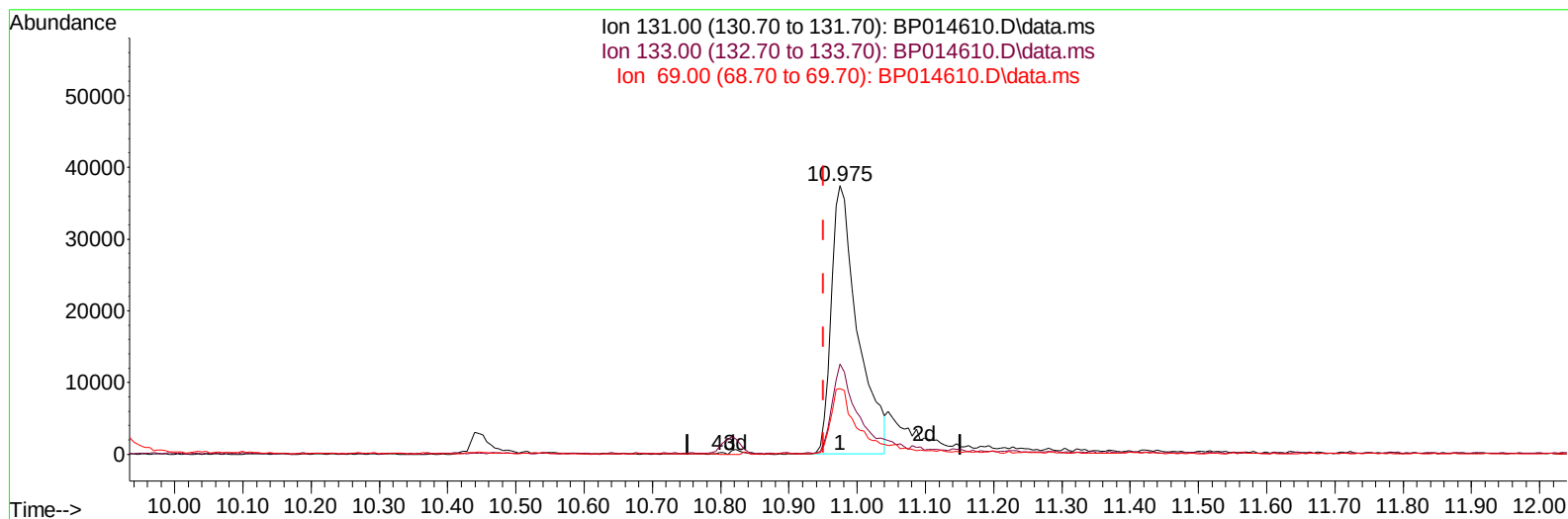
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP040723\
Data File : BP014610.D
Acq On : 07 Apr 2023 18:25
Operator : CG/JU
Sample : PB151979BL
Misc :
ALS Vial : 3 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
SBLK979

Manual IntegrationsAPPROVED

Reviewed By :Christian Giraldo 04/10/2023
Supervised By :Jagrut Upadhyay 04/10/2023

Quant Time: Apr 08 01:50:01 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP032823.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri Apr 07 00:45:34 2023
Response via : Initial Calibration



TIC: BP014610.D\data.ms

(31) 4-Chloroaniline-d4 (S)

10.975min (+ 0.024) 9.15 ng/u1

response 100202

Ion	Exp%	Act%
131.00	100.00	100.00
133.00	32.20	33.57
69.00	21.30	24.50
0.00	0.00	0.00

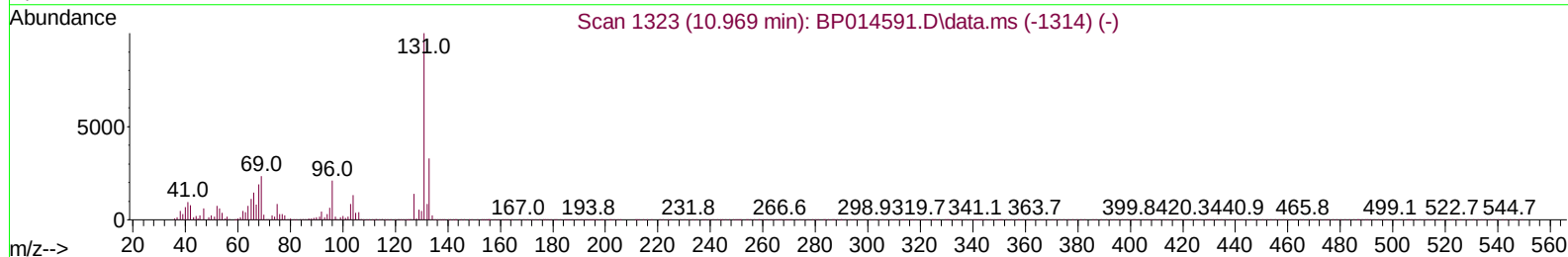
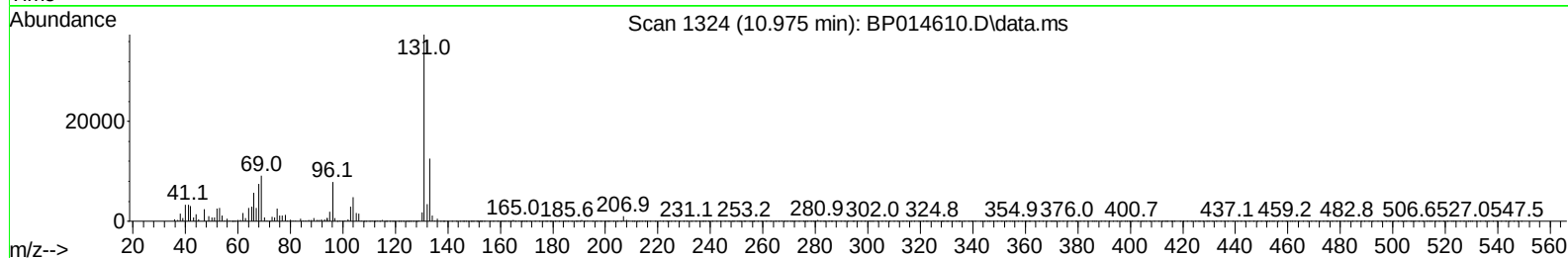
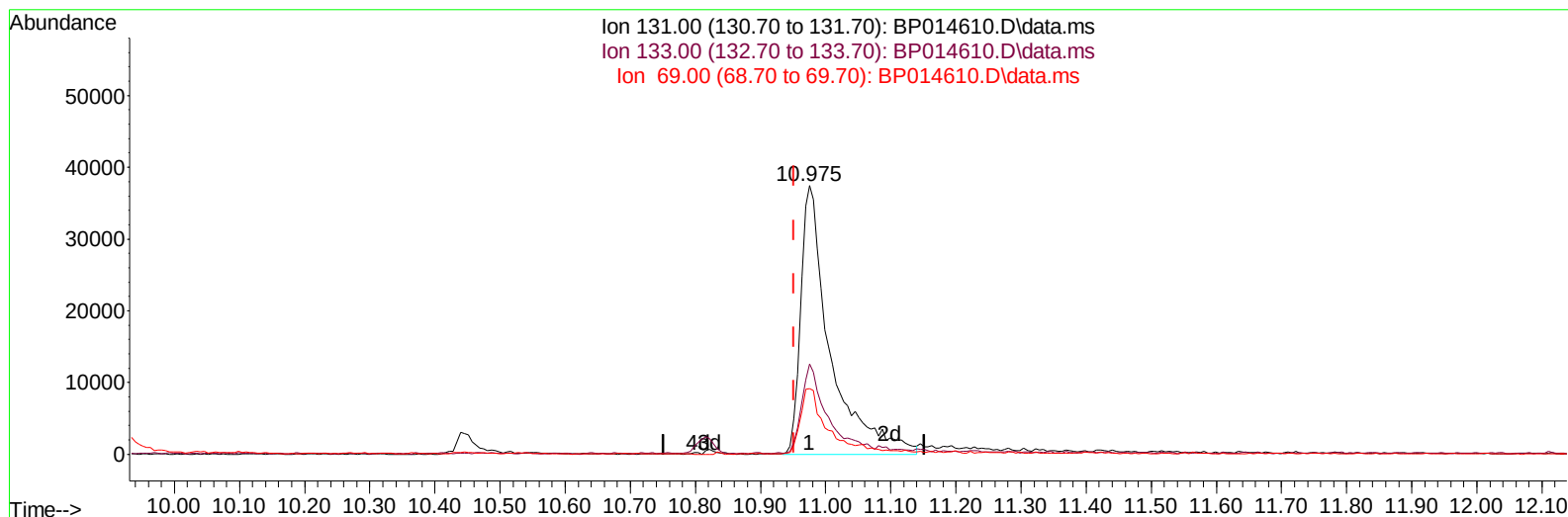
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(31) 4-Chloroaniline-d4 (S)

10.975min (+ 0.024) 10.69 ng/ul m

response 117066

Ion	Exp%	Act%
131.00	100.00	100.00
133.00	32.20	33.57
69.00	21.30	24.50
0.00	0.00	0.00

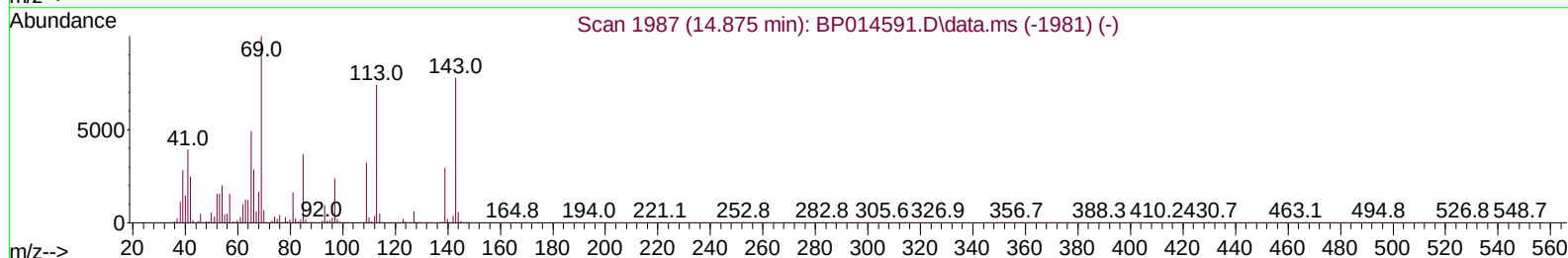
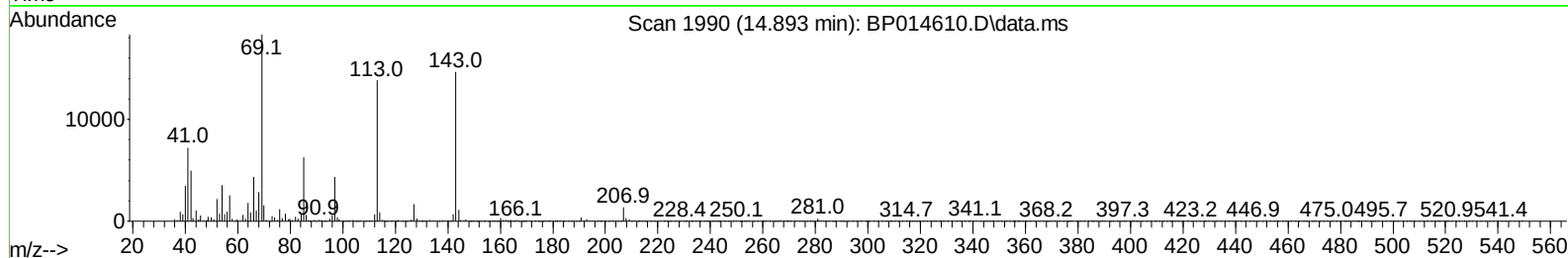
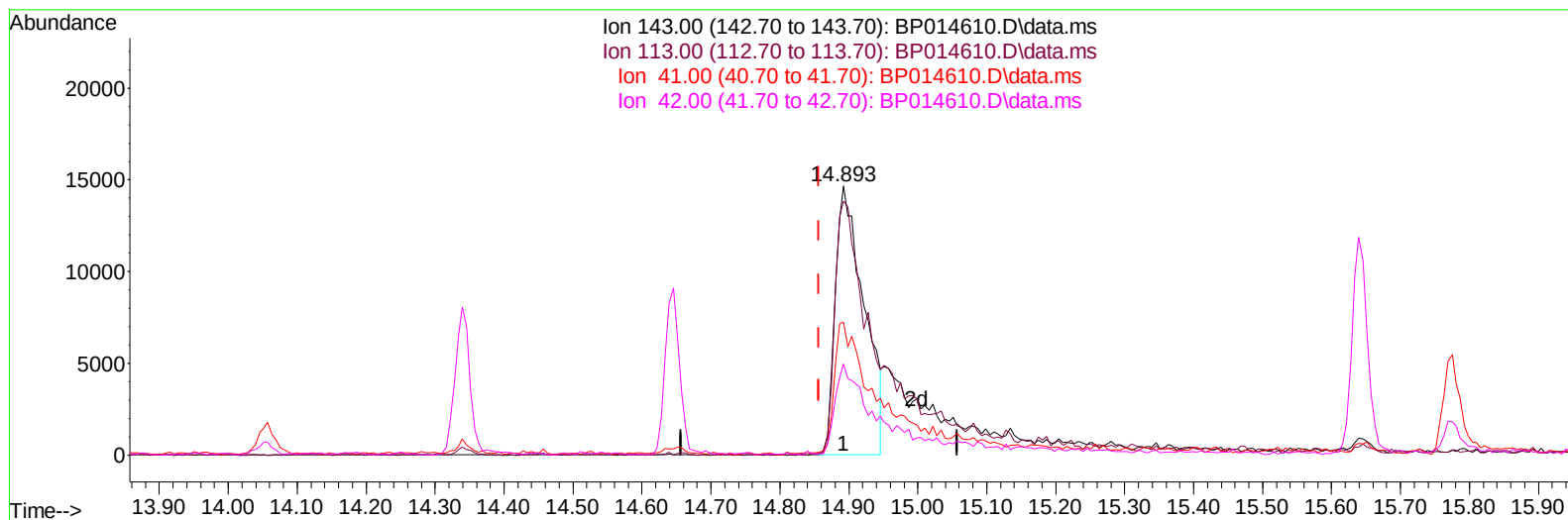
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TIC: BP014610.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.893min (+ 0.035) 10.80 ng/ul

response 42414

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	96.80	94.32
41.00	44.20	49.31
42.00	31.50	33.85

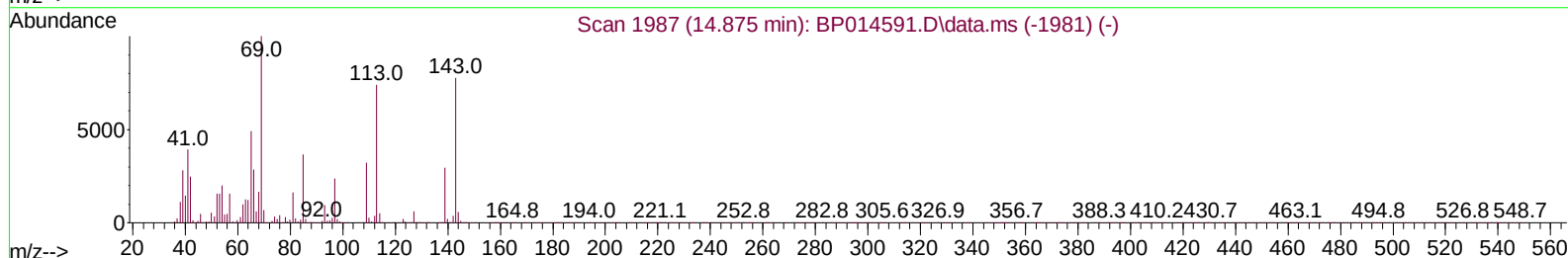
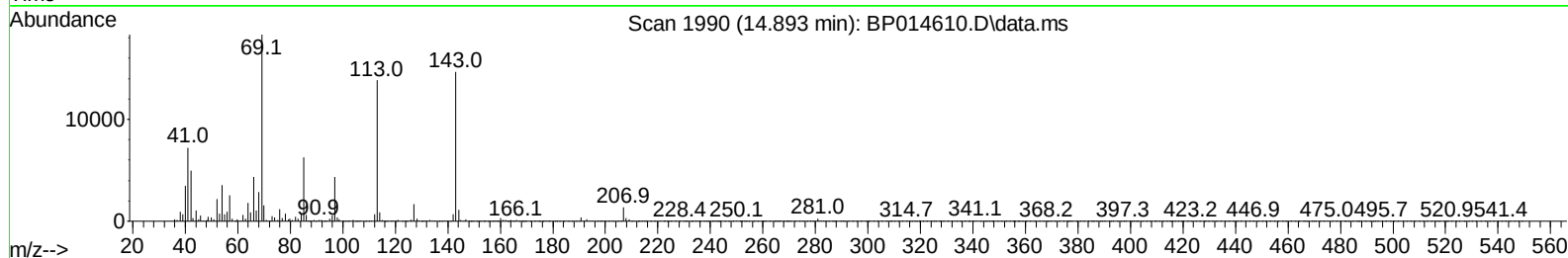
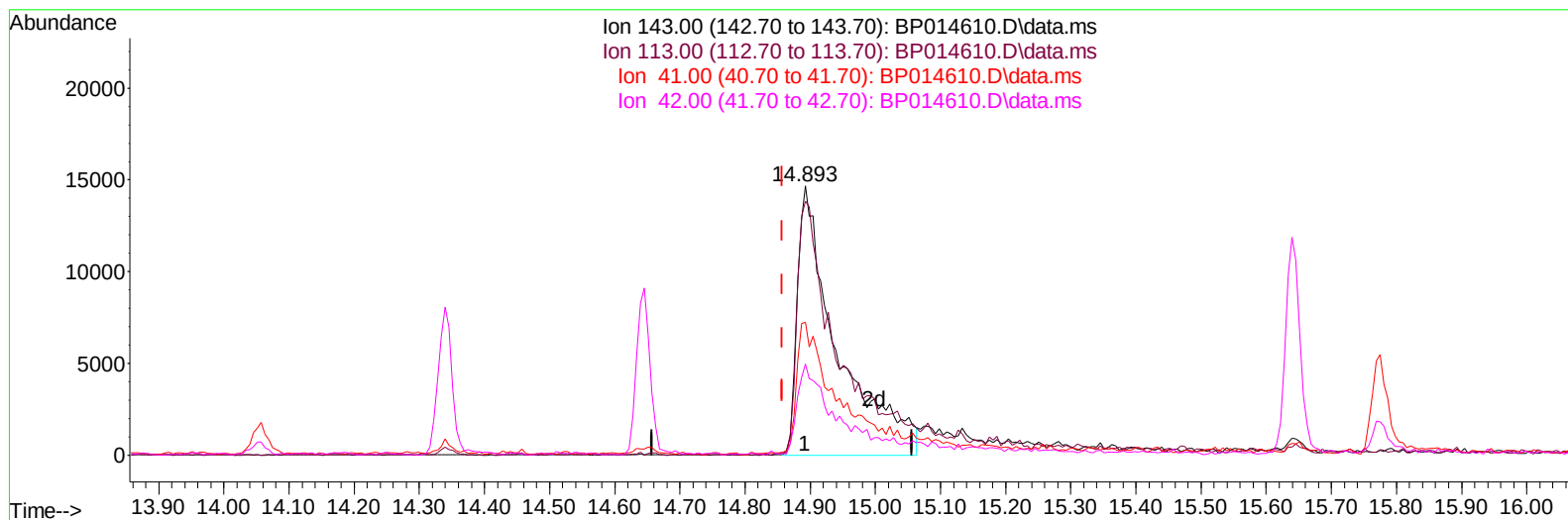
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Operator : CG/JU
Sample : PB151979BL
Misc :
ALS Vial : 3 Sample Multiplier: 1

Instrument :
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TIC: BP014610.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.893min (+ 0.035) 15.95 ng/ul m

response 62630

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	96.80	94.32
41.00	44.20	49.31
42.00	31.50	33.85

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.993	152	120989	20.000	ng/ul	0.01
20) Naphthalene-d8	10.810	136	493738	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.645	164	320553	20.000	ng/ul	0.01
64) Phenanthrene-d10	17.410	188	714211	20.000	ng/ul	0.01
79) Chrysene-d12	21.498	240	692440	20.000	ng/ul	0.01
88) Perylene-d12	24.004	264	817170	20.000	ng/ul	0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.370	96	17452	5.594	ng/uL	0.00
4) Pyridine-d5	3.811	84	100248	12.538	ng/ul	0.02
7) Phenol-d5	7.158	99	258767	23.191	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.322	67	187273	26.159	ng/ul	0.01
11) 2-Chlorophenol-d4	7.522	132	205064	25.427	ng/ul	0.01
15) 4-Methylphenol-d8	8.716	113	200382	22.263	ng/ul	0.01
21) Nitrobenzene-d5	9.169	128	94474	23.702	ng/ul	0.01
24) 2-Nitrophenol-d4	9.899	143	108556	23.785	ng/ul	0.01
28) 2,4-Dichlorophenol-d3	10.446	165	174544	20.999	ng/ul	0.02
31) 4-Chloroaniline-d4	10.975	131	117066m	10.692	ng/ul	0.02
46) Dimethylphthalate-d6	14.057	166	648885	24.966	ng/ul	0.02
49) Acenaphthylene-d8	14.340	160	714992	24.792	ng/ul	0.01
54) 4-Nitrophenol-d4	14.893	143	62630m	15.950	ng/ul	0.04
60) Fluorene-d10	15.640	176	547139	25.243	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.775	200	96265	18.318	ng/ul	0.01
73) Anthracene-d10	17.510	188	874102	25.289	ng/ul	0.01
81) Pyrene-d10	19.739	212	1061177	22.816	ng/ul	0.01
92) Benzo(a)pyrene-d12	23.839	264	1120665	25.812	ng/ul	0.02

Target Compounds Qvalue

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

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