

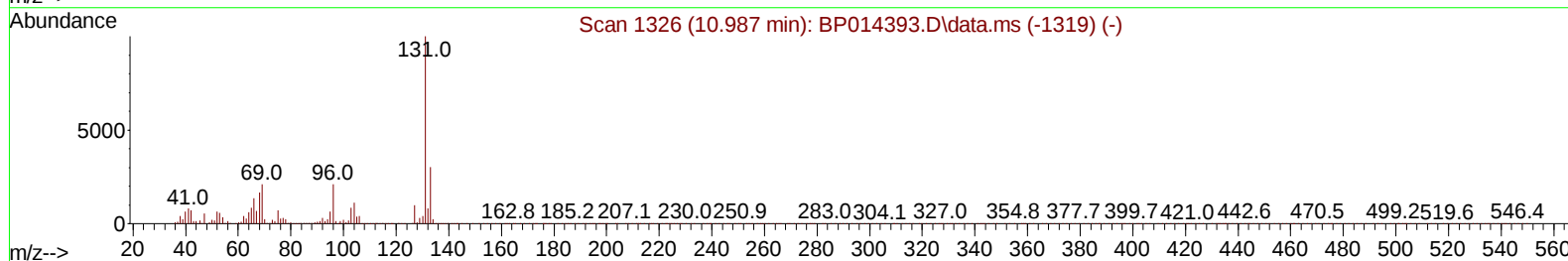
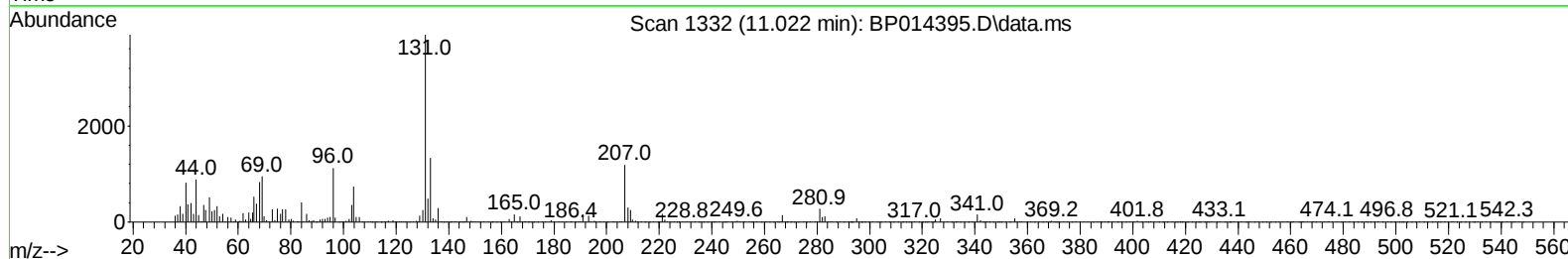
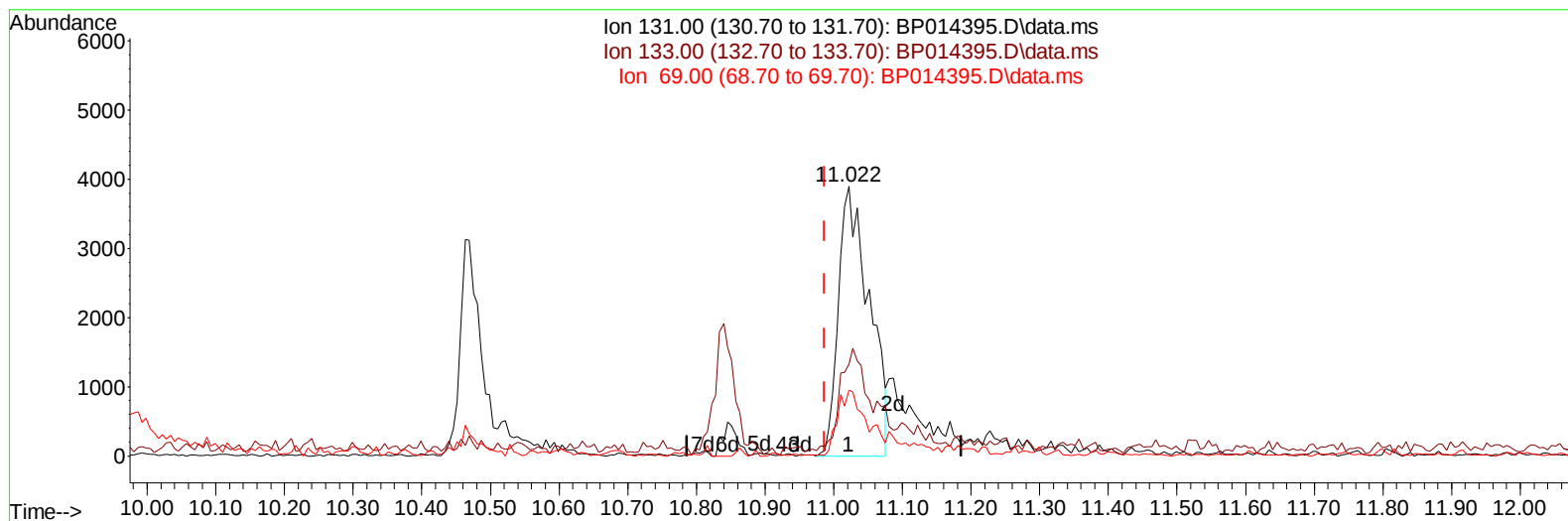
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP033023\
Data File : BP014395.D
Acq On : 30 Mar 2023 11:12
Operator : CG/JU
Sample : PB151778BL
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
SBLK778

Manual IntegrationsAPPROVED

Reviewed By :Christian Giraldo 03/31/2023
Supervised By :Jagrut Upadhyay 03/31/2023

Quant Time: Mar 30 22:52:16 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP032823.M
Quant Title : SVOA CALIBRATION
QLast Update : Thu Mar 30 12:32:14 2023
Response via : Initial Calibration



TIC: BP014395.D\data.ms

(31) 4-Chloroaniline-d4 (S)

11.022min (+ 0.035) 1.34 ng/u1

response 12029

Ion	Exp%	Act%
131.00	100.00	100.00
133.00	32.20	34.17
69.00	21.30	24.44
0.00	0.00	0.00

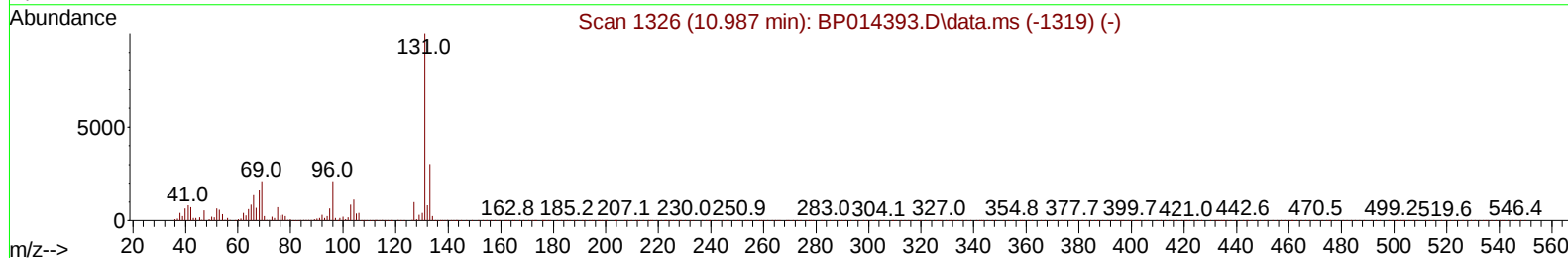
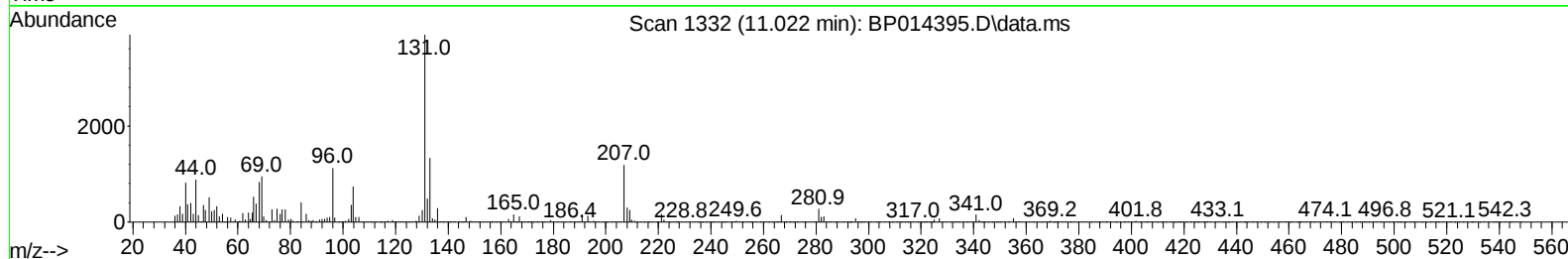
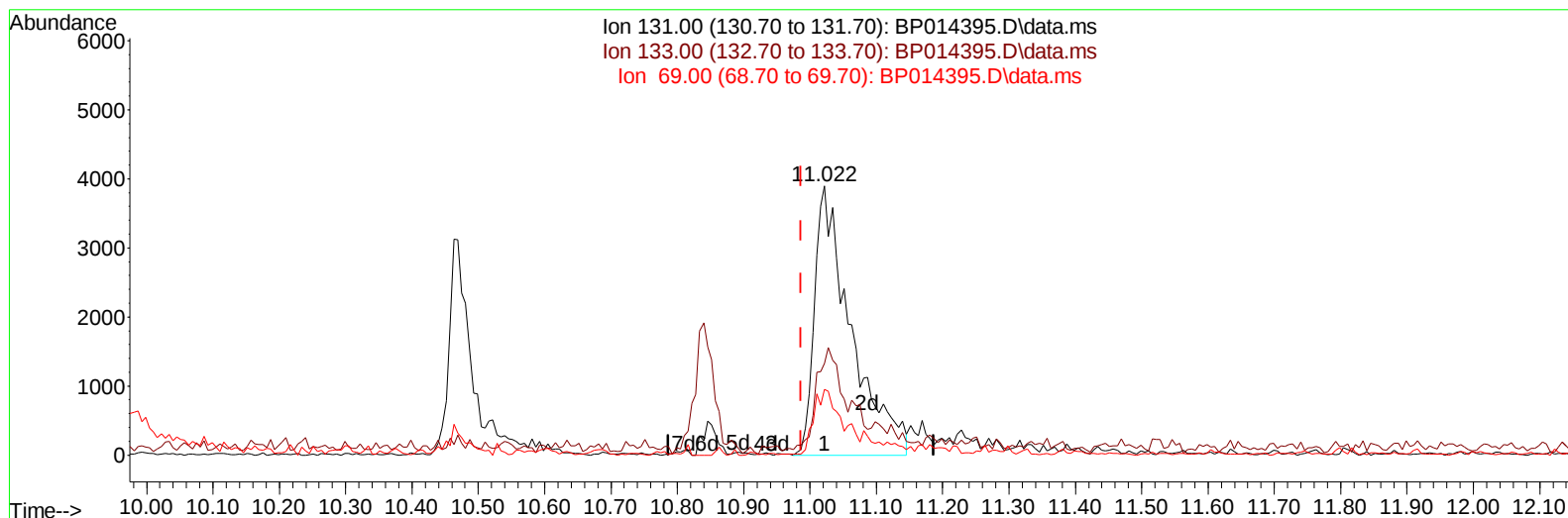
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(31) 4-Chloroaniline-d4 (S)

11.022min (+ 0.035) 1.65 ng/ul m

response 14834

Ion	Exp%	Act%
131.00	100.00	100.00
133.00	32.20	34.17
69.00	21.30	24.44
0.00	0.00	0.00

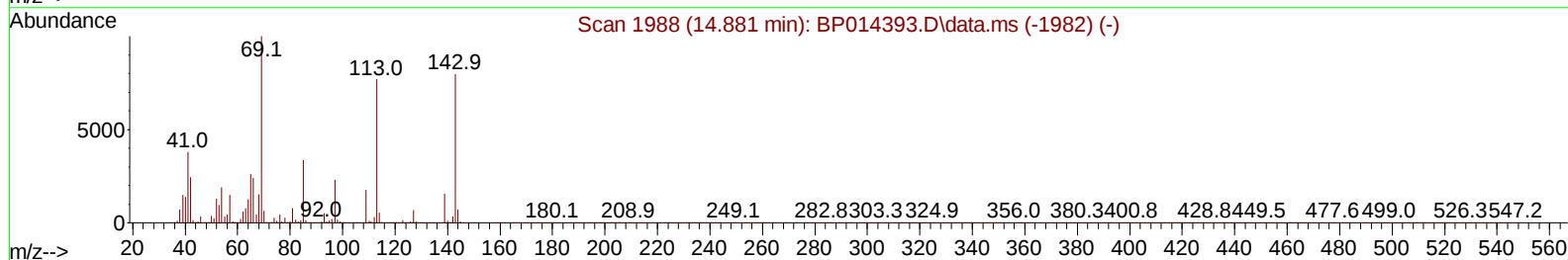
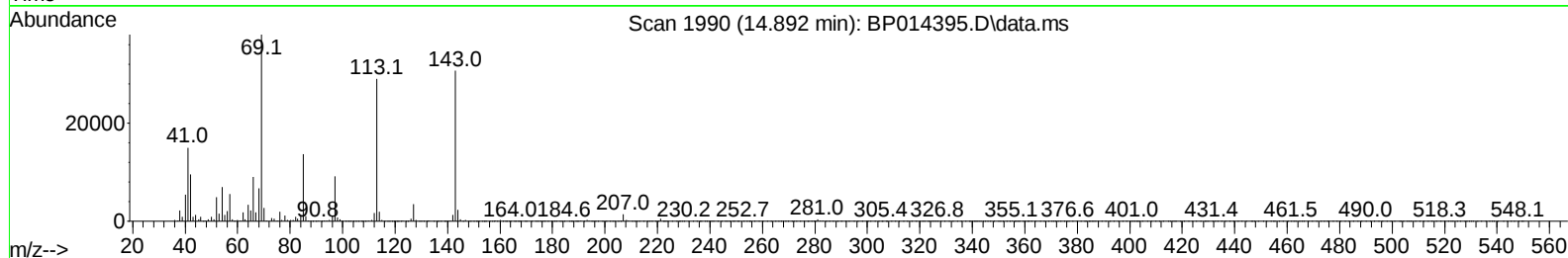
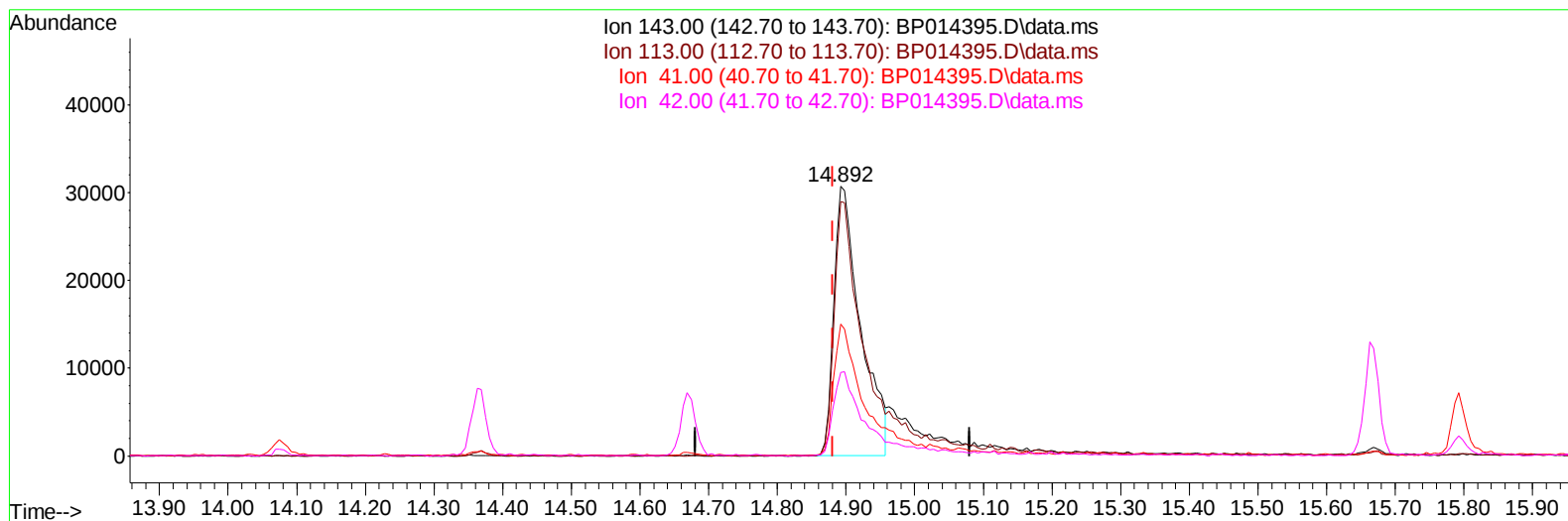
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TIC: BP014395.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.892min (+ 0.012) 23.95 ng/u1

response 82388

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	96.80	94.48
41.00	44.20	48.89
42.00	31.50	31.00

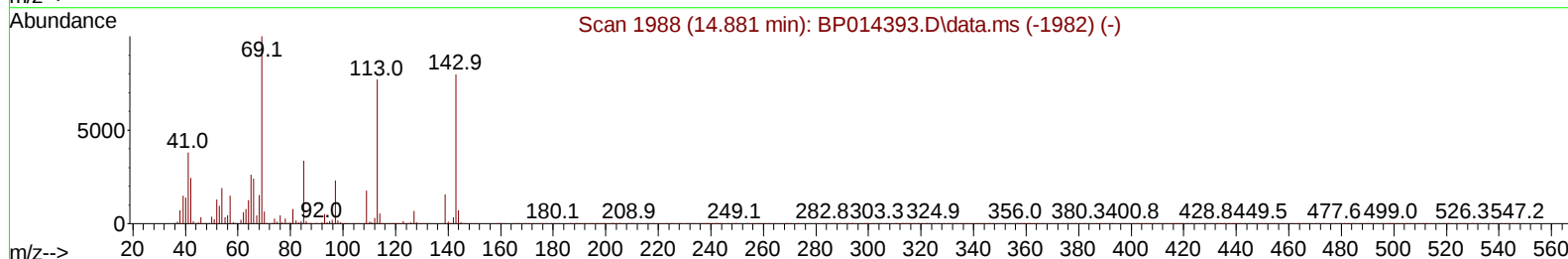
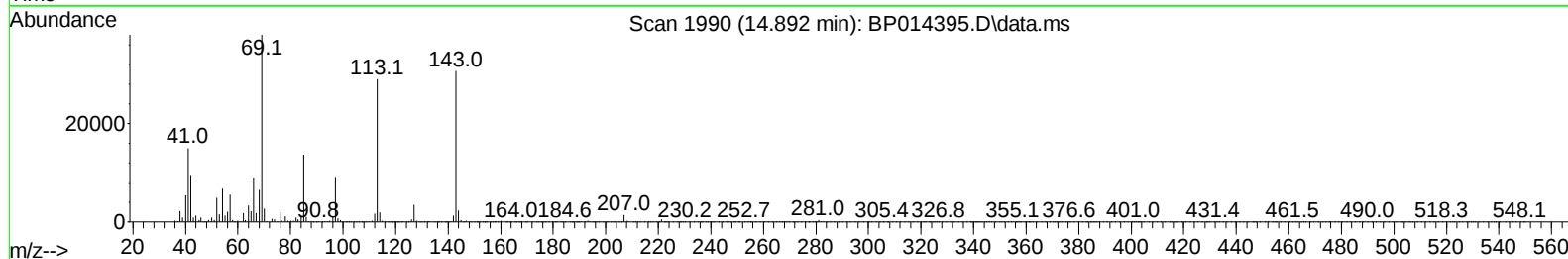
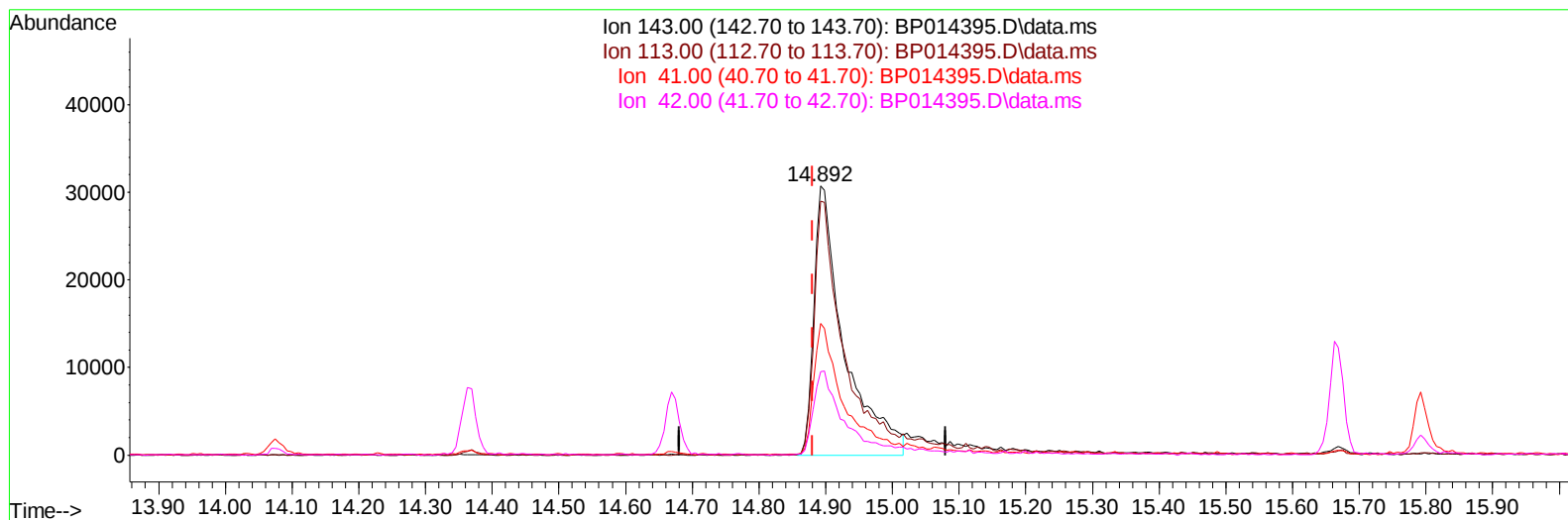
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Operator : CG/JU
Sample : PB151778BL
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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TIC: BP014395.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.892min (+ 0.012) 27.98 ng/ul m

response 96259

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	96.80	94.48
41.00	44.20	48.89
42.00	31.50	31.00

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.016	152	93398	20.000	ng/ul	0.00
20) Naphthalene-d8	10.840	136	406059	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.669	164	280832	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.433	188	671954	20.000	ng/ul	0.00
79) Chrysene-d12	21.521	240	705010	20.000	ng/ul	0.00
88) Perylene-d12	24.039	264	765944	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.393	96	15589	6.473	ng/uL	0.00
4) Pyridine-d5	3.828	84	86479	14.011	ng/ul	0.00
7) Phenol-d5	7.175	99	267503	31.056	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.340	67	180345	32.633	ng/ul	0.00
11) 2-Chlorophenol-d4	7.546	132	202639	32.549	ng/ul	0.00
15) 4-Methylphenol-d8	8.734	113	219588	31.604	ng/ul	0.00
21) Nitrobenzene-d5	9.193	128	99794	30.443	ng/ul	0.00
24) 2-Nitrophenol-d4	9.922	143	112538	29.981	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.469	165	187959	27.495	ng/ul	0.00
31) 4-Chloroaniline-d4	11.022	131	14834m	1.647	ng/ul	0.04
46) Dimethylphthalate-d6	14.075	166	756672	33.231	ng/ul	0.00
49) Acenaphthylene-d8	14.363	160	800965	31.701	ng/ul	0.00
54) 4-Nitrophenol-d4	14.892	143	96259m	27.982	ng/ul	0.01
60) Fluorene-d10	15.669	176	645682	34.003	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.792	200	123003	24.878	ng/ul	0.00
73) Anthracene-d10	17.533	188	1097838	33.759	ng/ul	0.00
81) Pyrene-d10	19.763	212	1333338	28.156	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.880	264	1365194	33.548	ng/ul	0.00

Target Compounds Qvalue

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

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