

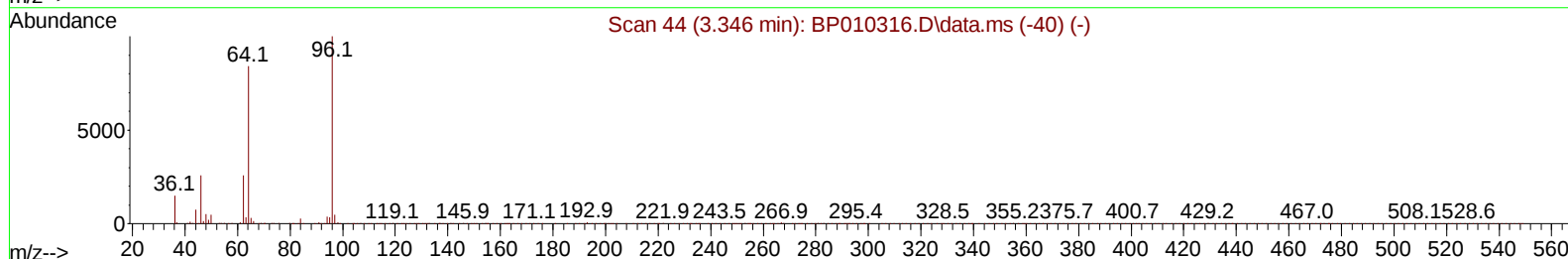
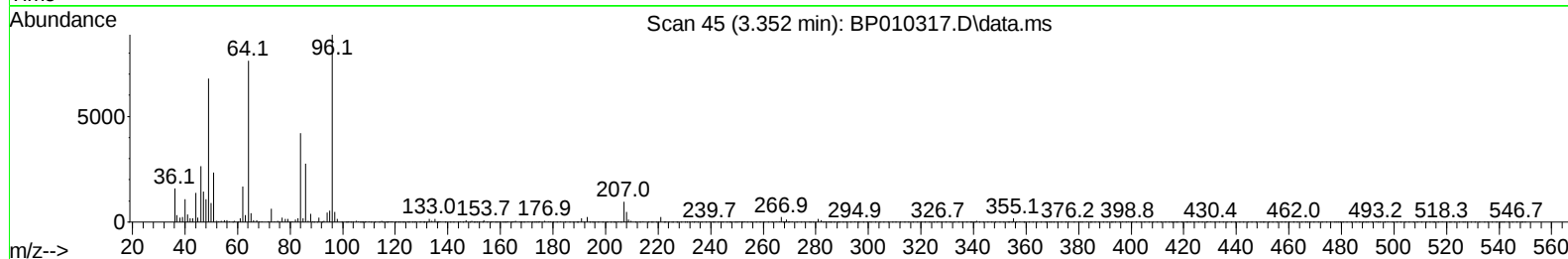
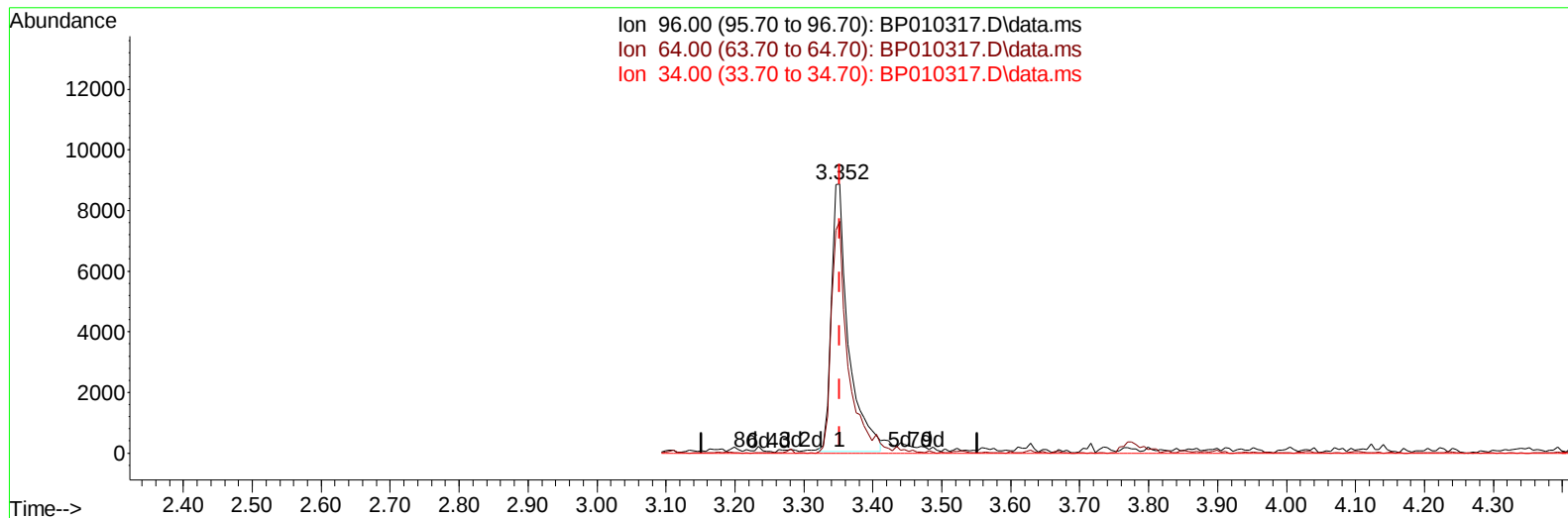
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051122\
 Data File : BP010317.D
 Acq On : 12 May 2022 10:42
 Operator : CG/JU
 Sample : PB144712BL
 Misc :
 ALS Vial : 41 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK712

Manual IntegrationsAPPROVED

Quant Time: May 12 23:59:53 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 05/16/2022
 Supervised By :Yogesh Patel 05/17/2022



TIC: BP010317.D\data.ms

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

3.352min (-0.000) 5.78 ng/uL

response 15125

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	86.14#
34.00	0.00	0.00
0.00	0.00	0.00

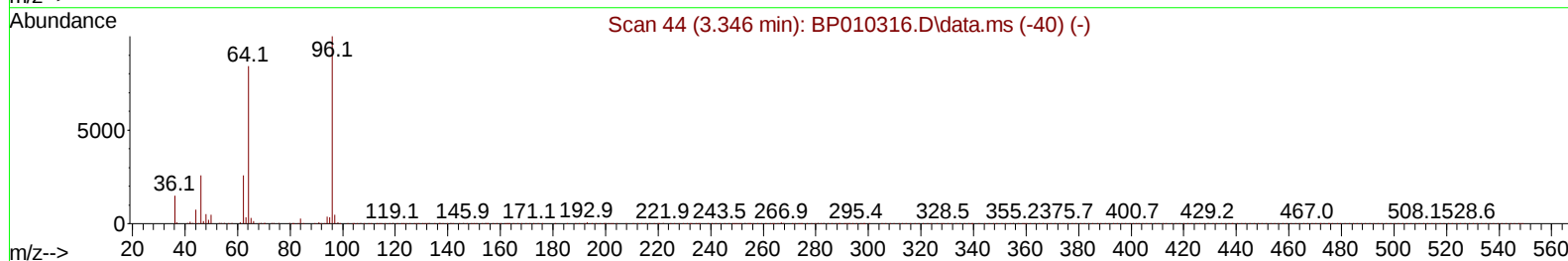
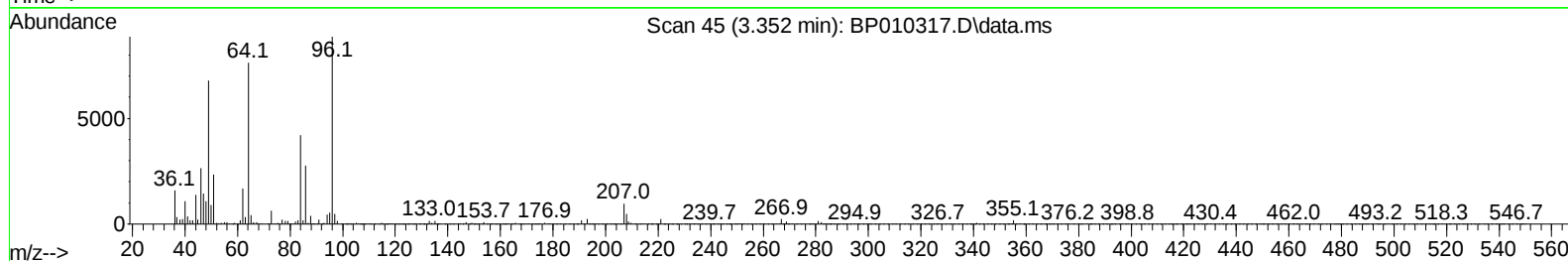
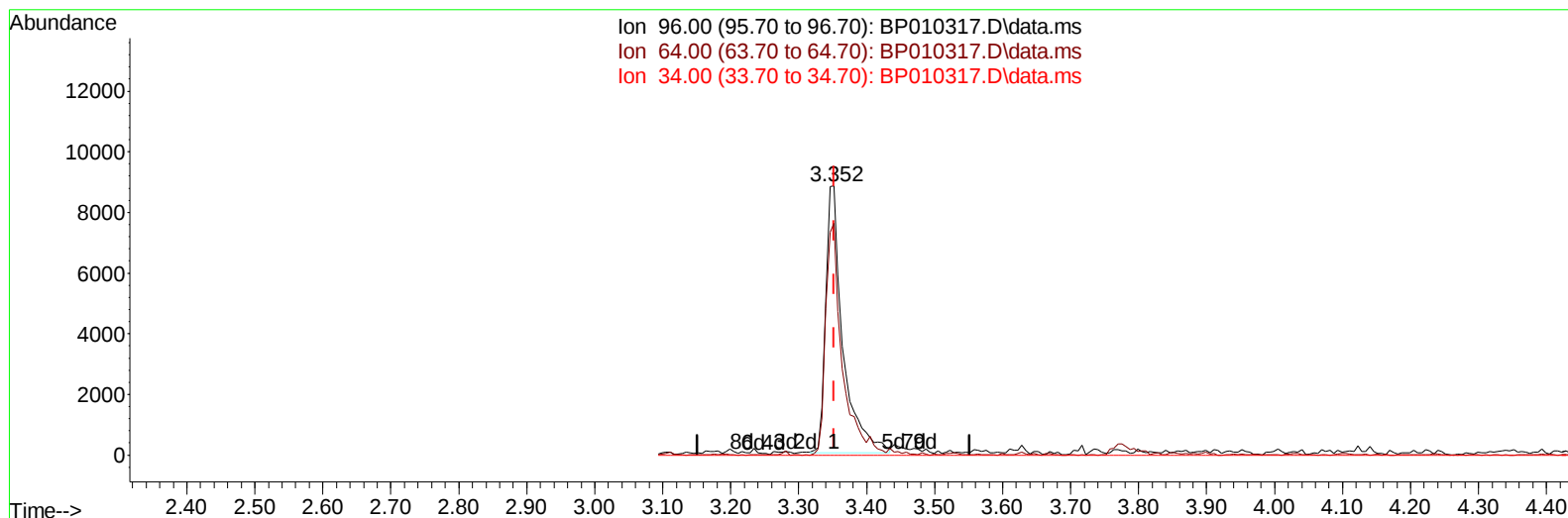
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TIC: BP010317.D\data.ms

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

3.352min (-0.000) 5.90 ng/uL m

response 15450

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	86.14#
34.00	0.00	0.00
0.00	0.00	0.00

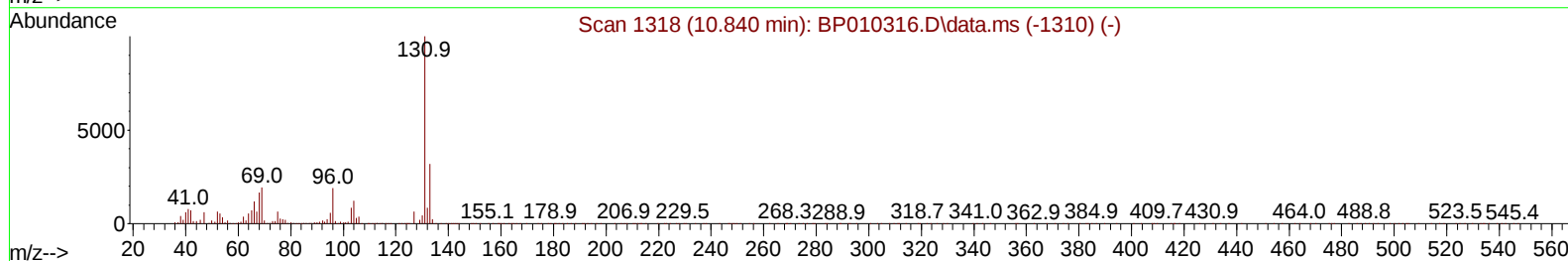
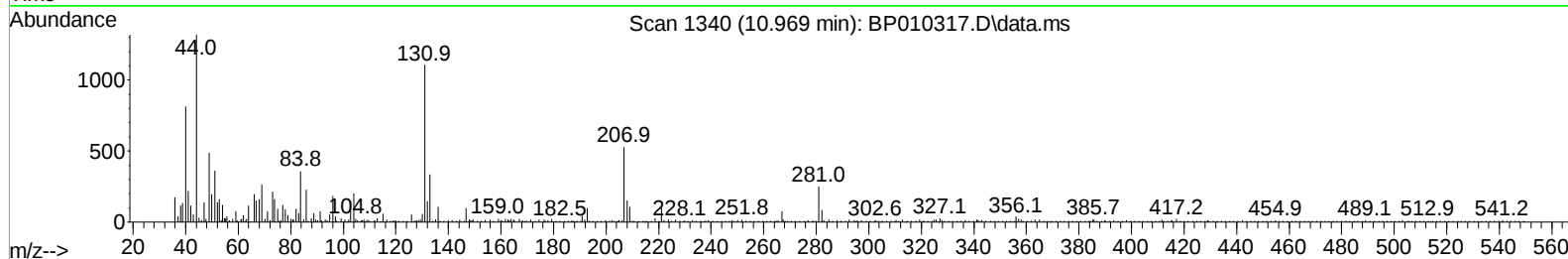
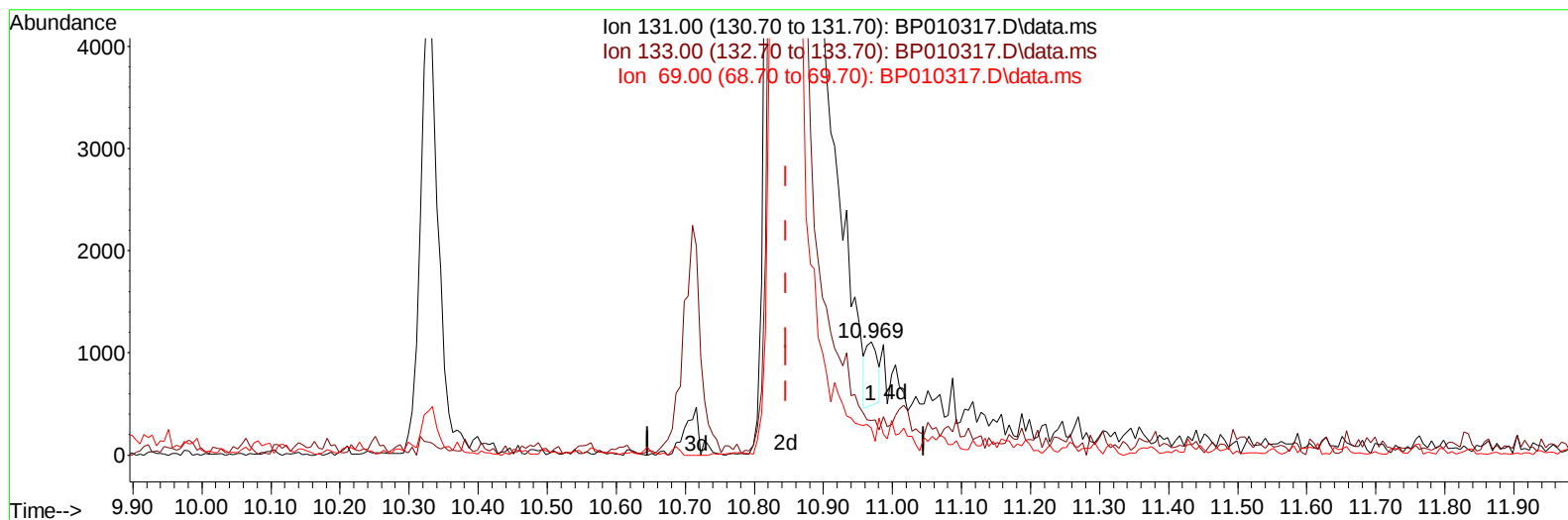
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TIC: BP010317.D\data.ms

(31) 4-Chloroaniline-d4 (S)

10.969min (+ 0.124) 0.06 ng/ul

response 744

Ion	Exp%	Act%
131.00	100.00	100.00
133.00	29.20	30.44
69.00	26.10	23.94
0.00	0.00	0.00

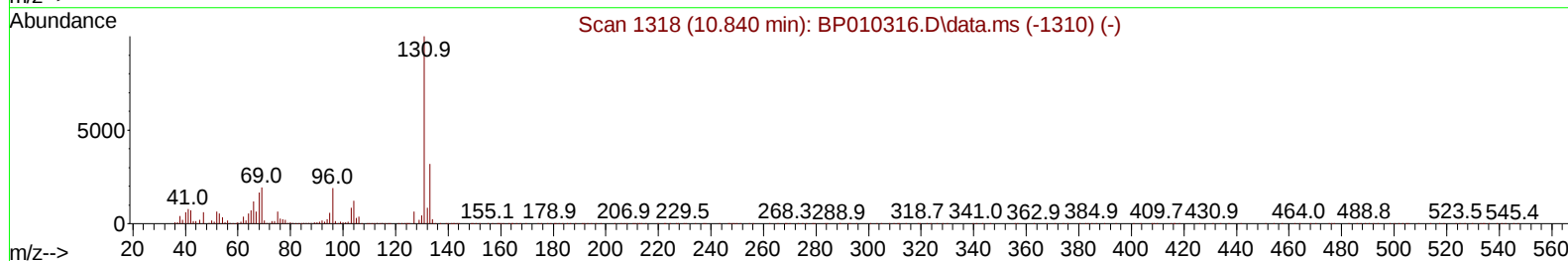
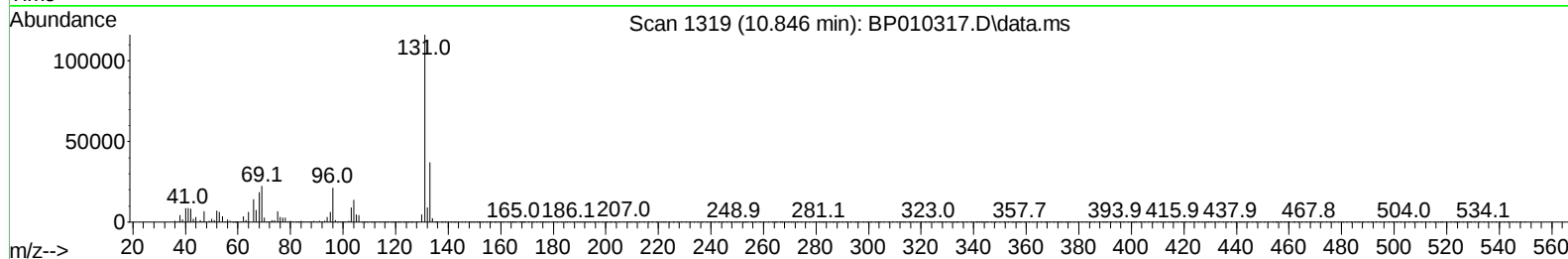
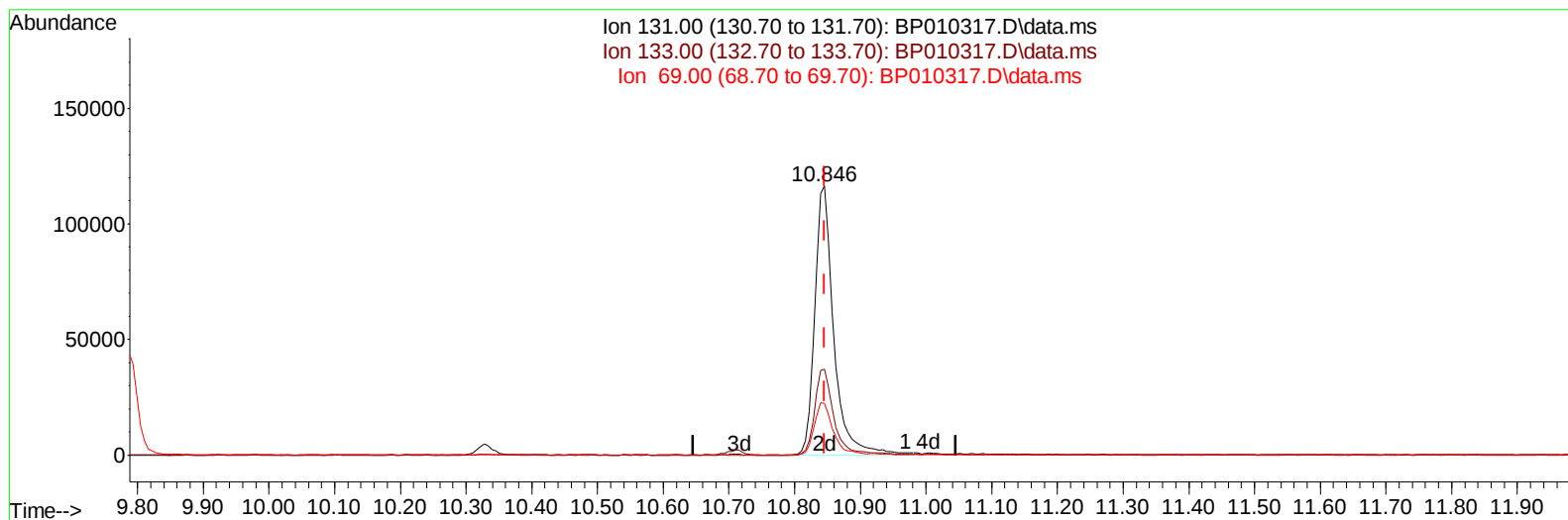
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Misc :
ALS Vial : 41 Sample Multiplier: 1

Instrument :
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Manual IntegrationsAPPROVED

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TIC: BP010317.D\data.ms

(31) 4-Chloroaniline-d4 (S)

10.846min (-0.000) 19.23 ng/ul m

response 235924

Ion	Exp%	Act%
131.00	100.00	100.00
133.00	29.20	32.02
69.00	26.10	19.27#
0.00	0.00	0.00

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Compound	R.T.	Q	Ion	Response	Conc	Units	Dev(Min)

Internal Standards							
1) 1,4-Dichlorobenzene-d4	7.905	152		111182	20.000	ng/ul	0.00
20) Naphthalene-d8	10.704	136		499528	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.534	164		358109	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.286	188		805612	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.369	240		814745	20.000	ng/ul	# 0.00
88) Perylene-d12	23.798	264		810710	20.000	ng/ul	-0.02
System Monitoring Compounds							
3) 1,4-Dioxane-d8	3.352	96		15450m	5.903	ng/uL	0.00
4) Pyridine-d5	3.769	84		179621	24.663	ng/ul	0.00
7) Phenol-d5	7.069	99		273383	26.474	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.234	67		194844	30.501	ng/ul	0.00
11) 2-Chlorophenol-d4	7.434	132		219265	28.831	ng/ul	0.00
15) 4-Methylphenol-d8	8.610	113		233408	27.493	ng/ul	0.00
21) Nitrobenzene-d5	9.063	128		117283	29.282	ng/ul	0.00
24) 2-Nitrophenol-d4	9.787	143		127262	27.867	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.328	165		217502	26.167	ng/ul	0.00
31) 4-Chloroaniline-d4	10.846	131		235924m	19.226	ng/ul	0.00
46) Dimethylphthalate-d6	13.945	166		858497	30.092	ng/ul	0.00
49) Acenaphthylene-d8	14.228	160		936987	27.547	ng/ul	0.00
54) 4-Nitrophenol-d4	14.734	143		121174	22.143	ng/ul	0.00
60) Fluorene-d10	15.528	176		719419	29.874	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.651	200		122125	22.137	ng/ul	0.00
73) Anthracene-d10	17.386	188		1198522	30.380	ng/ul	0.00
81) Pyrene-d10	19.616	212		1457079	32.363	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.645	264		1361863	31.499	ng/ul	-0.01

Target Compounds Qvalue

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

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