

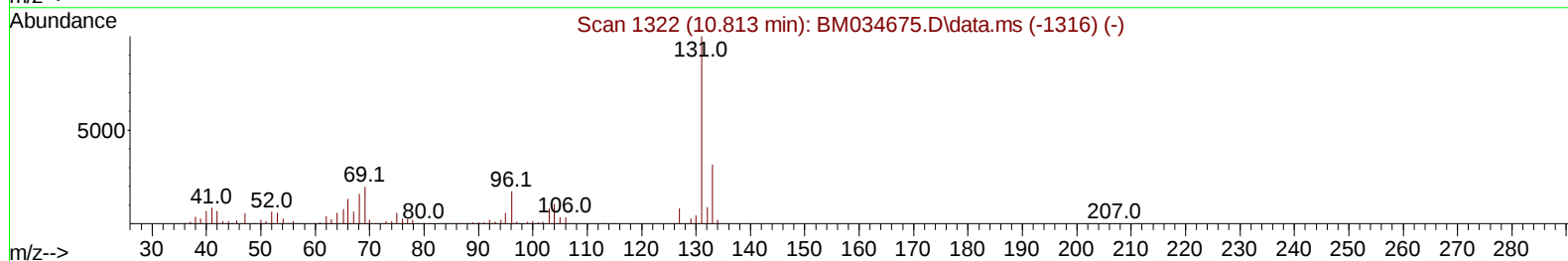
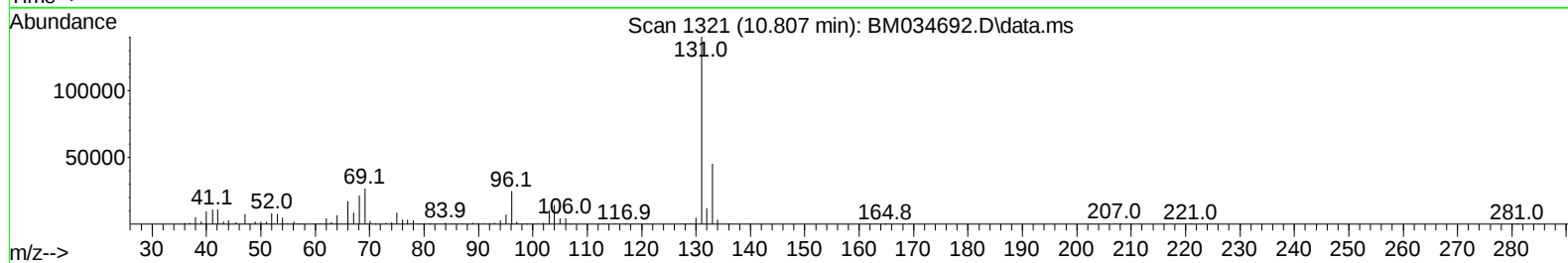
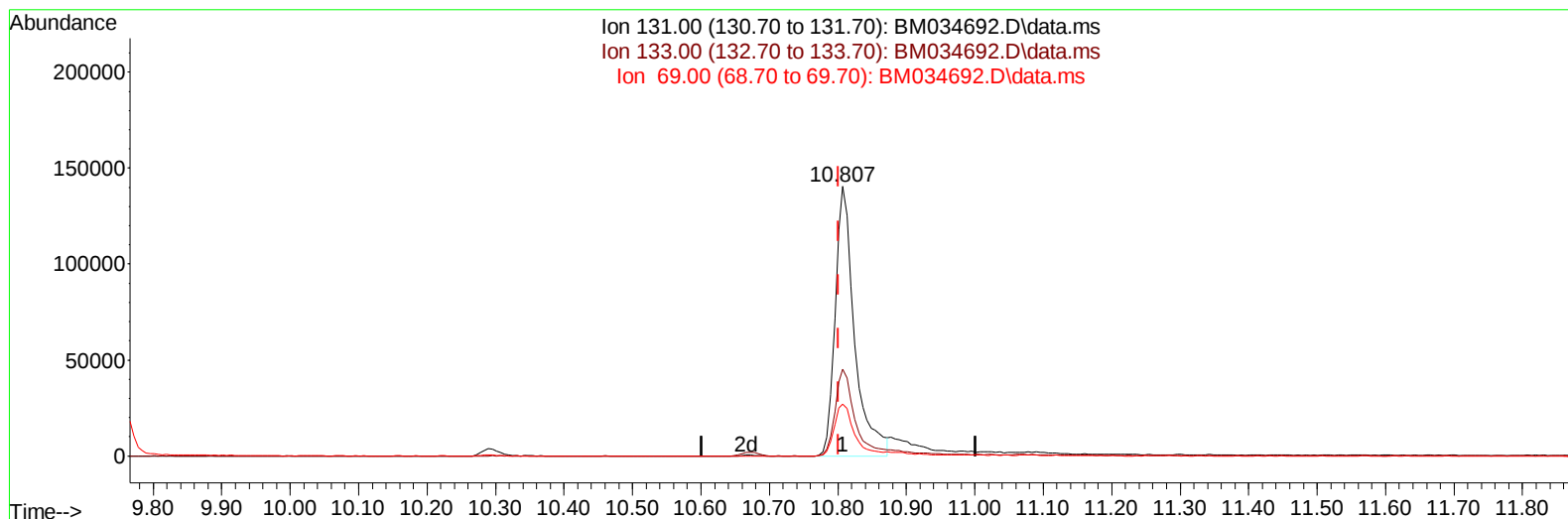
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM041422\
Data File : BM034692.D
Acq On : 15 Apr 2022 06:39
Operator : CG/JU
Sample : PB144055BL
Misc :
ALS Vial : 35 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
SBLK055

Manual IntegrationsAPPROVED

Quant Time: Apr 18 06:29:40 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM040922.M
Quant Title : SVOA CALIBRATION
QLast Update : Mon Apr 18 05:39:41 2022
Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 04/22/2022
Supervised By :Yogesh Patel 04/23/2022



TIC: BM034692.D\data.ms

(31) 4-Chloroaniline-d4 (S)

10.807min (+ 0.006) 28.35 ng/u1

response 276751

Ion	Exp%	Act%
131.00	100.00	100.00
133.00	33.00	32.32
69.00	16.80	19.25
0.00	0.00	0.00

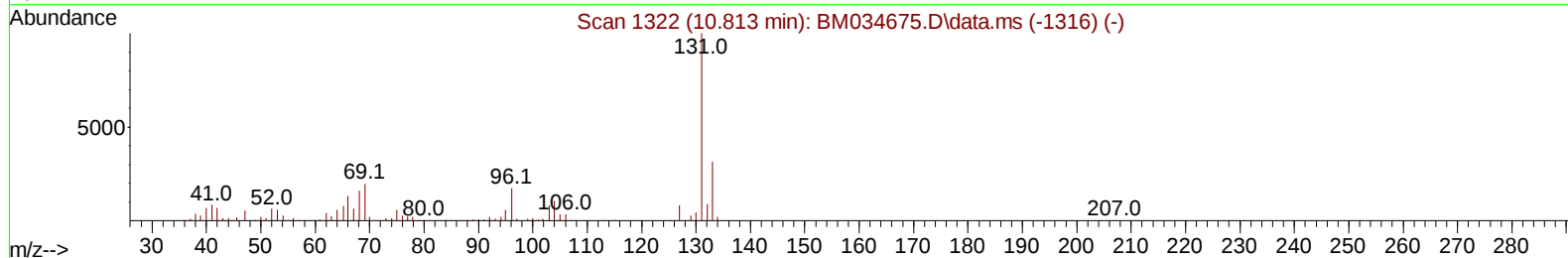
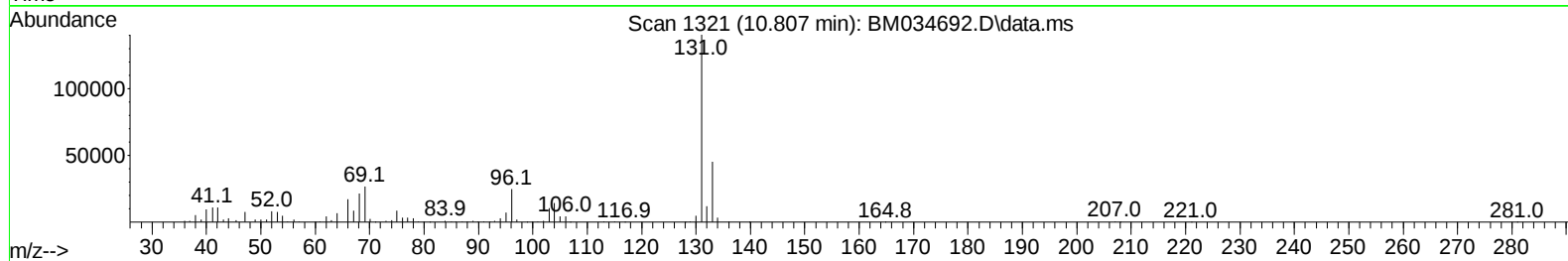
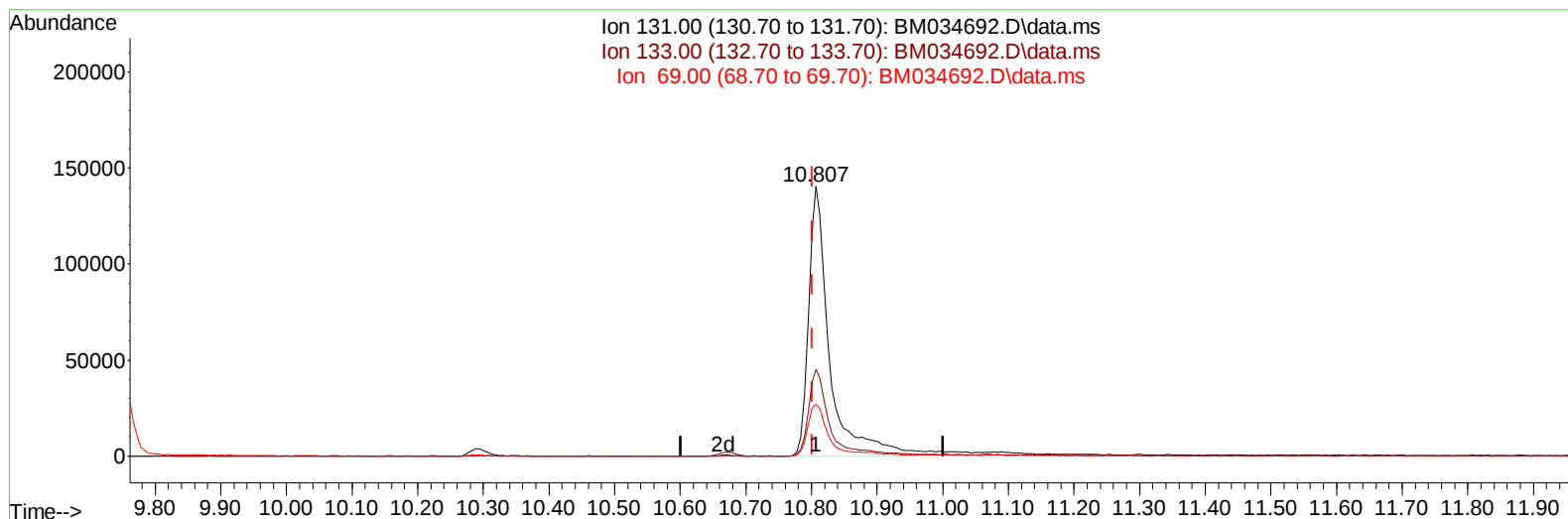
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TIC: BM034692.D\data.ms

(31) 4-Chloroaniline-d4 (S)

10.807min (+ 0.006) 31.33 ng/ul m

response 305876

Ion	Exp%	Act%
131.00	100.00	100.00
133.00	33.00	32.32
69.00	16.80	19.25
0.00	0.00	0.00

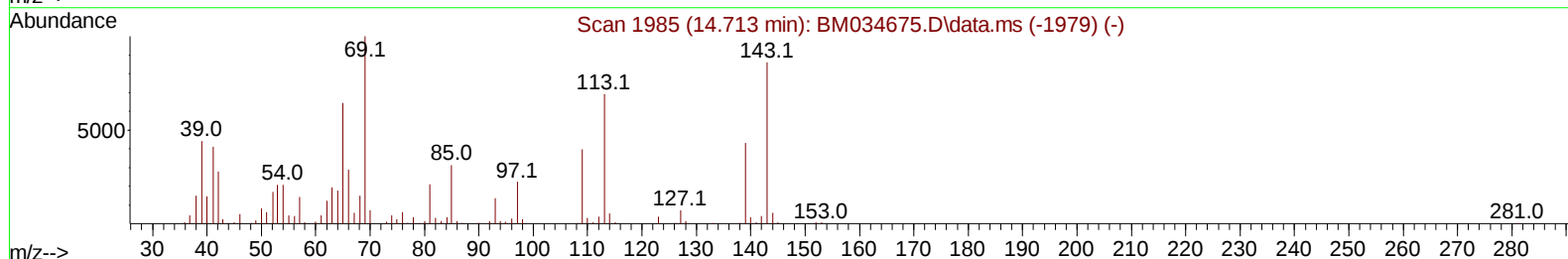
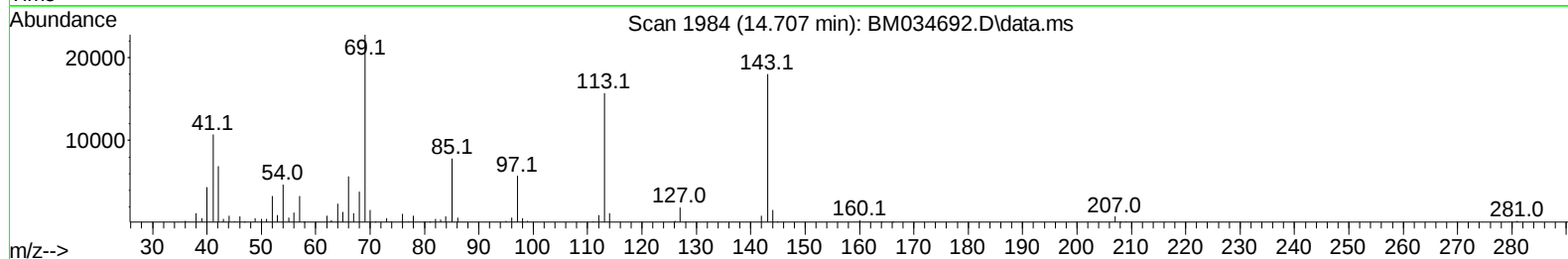
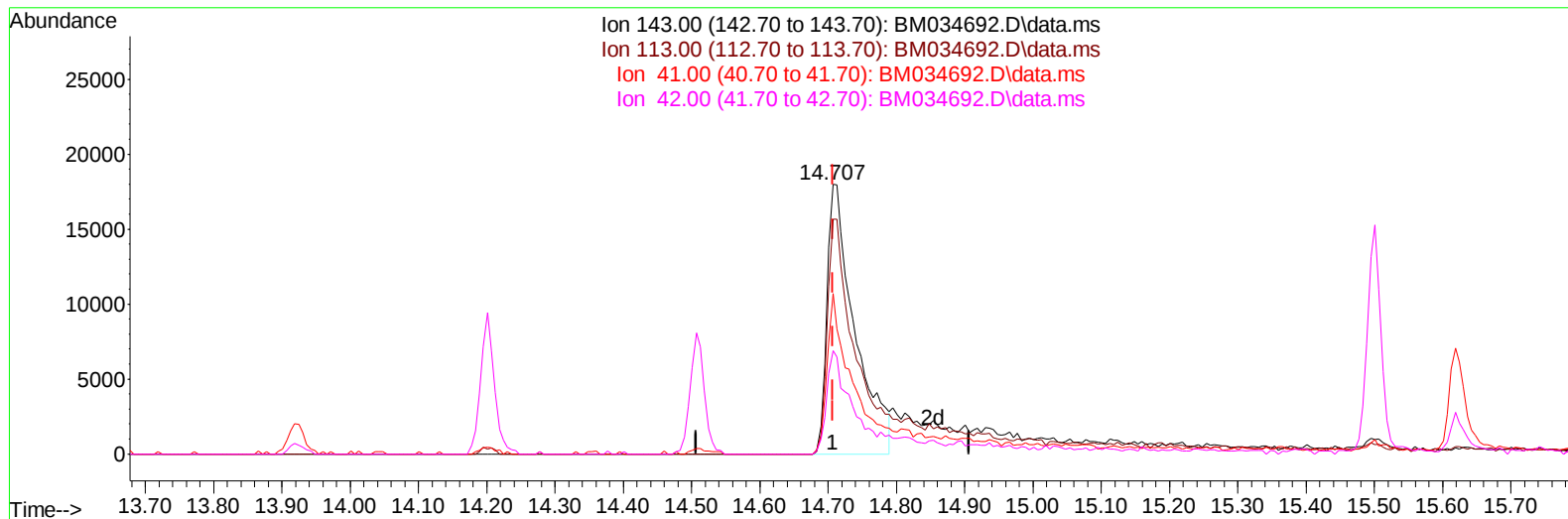
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Instrument :
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TIC: BM034692.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.707min (+ 0.000) 15.86 ng/u1

response 50973

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	77.80	87.13
41.00	31.90	59.49#
42.00	19.10	38.30#

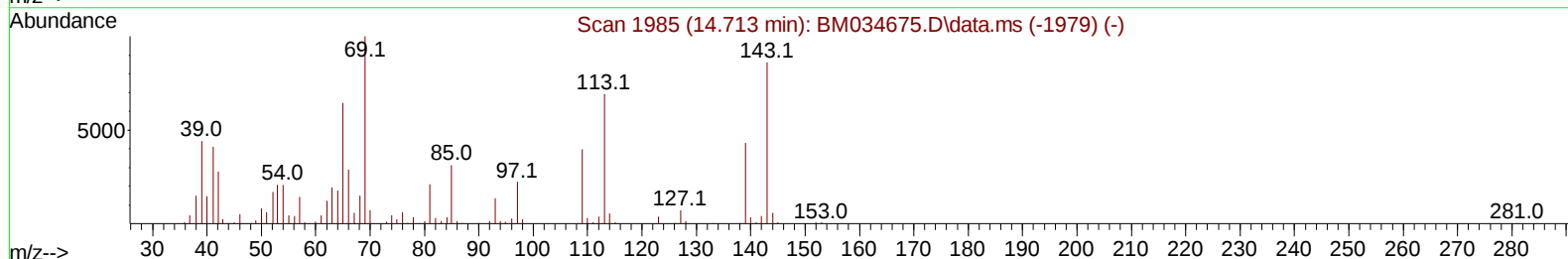
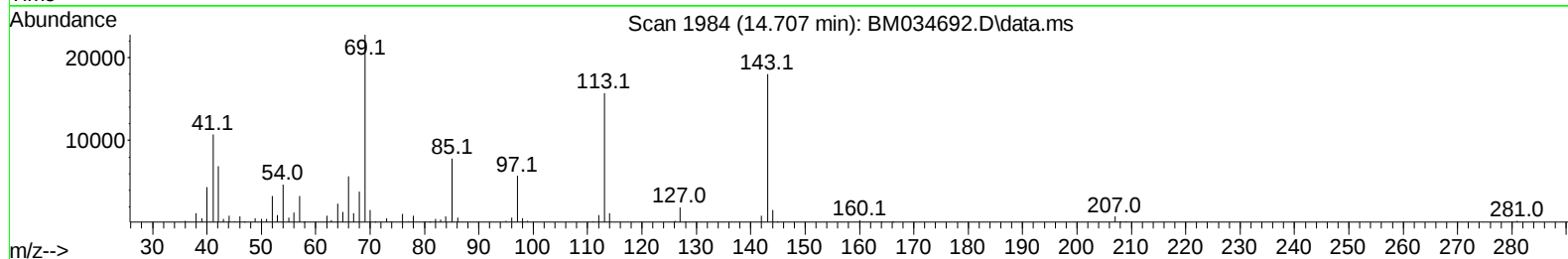
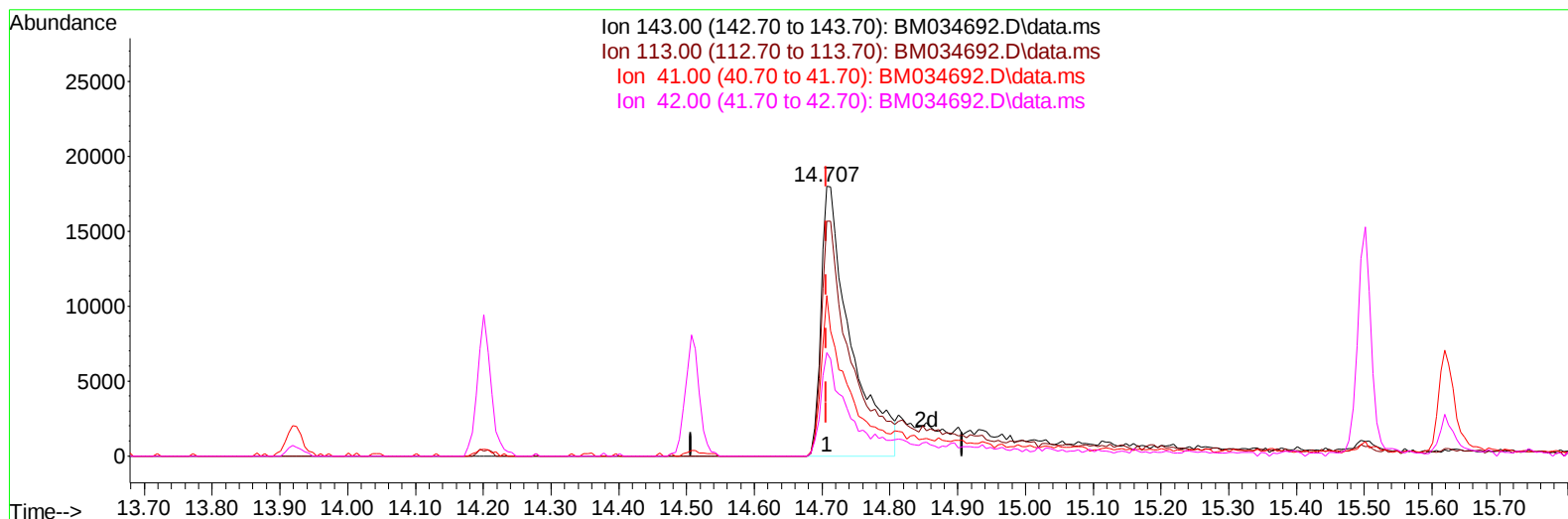
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Misc :
ALS Vial : 35 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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Supervised By :Yogesh Patel 04/23/2022



TIC: BM034692.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.707min (+ 0.000) 16.73 ng/ul m

response 53797

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	77.80	87.13
41.00	31.90	59.49#
42.00	19.10	38.30#

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.860	152	116132	20.000	ng/ul	0.00
20) Naphthalene-d8	10.672	136	491117	20.000	ng/ul	# 0.01
38) Acenaphthene-d10	14.507	164	291789	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.254	188	563763	20.000	ng/ul	# 0.01
79) Chrysene-d12	21.430	240	527031	20.000	ng/ul	# 0.00
88) Perylene-d12	23.800	264	569743	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.243	96	21684	6.658	ng/uL	0.01
4) Pyridine-d5	3.667	84	257502	32.763	ng/ul	0.00
7) Phenol-d5	7.025	99	305592	31.616	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.190	67	229432	35.621	ng/ul	0.01
11) 2-Chlorophenol-d4	7.390	132	266250	35.633	ng/ul	0.00
15) 4-Methylphenol-d8	8.572	113	253719	33.603	ng/ul	0.00
21) Nitrobenzene-d5	9.025	128	124811	34.579	ng/ul	0.01
24) 2-Nitrophenol-d4	9.748	143	131558	35.009	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.295	165	215963	30.520	ng/ul	0.01
31) 4-Chloroaniline-d4	10.807	131	305876m	31.332	ng/ul	0.00
46) Dimethylphthalate-d6	13.919	166	743384	35.258	ng/ul	0.00
49) Acenaphthylene-d8	14.201	160	885896	32.691	ng/ul	0.00
54) 4-Nitrophenol-d4	14.707	143	53797m	16.734	ng/ul	0.00
60) Fluorene-d10	15.501	176	621349	34.102	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.619	200	94921	27.580	ng/ul	0.00
73) Anthracene-d10	17.354	188	950977	35.381	ng/ul	0.01
81) Pyrene-d10	19.642	212	1096150	38.107	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.648	264	1001457	34.943	ng/ul	0.00

Target Compounds Qvalue

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

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