

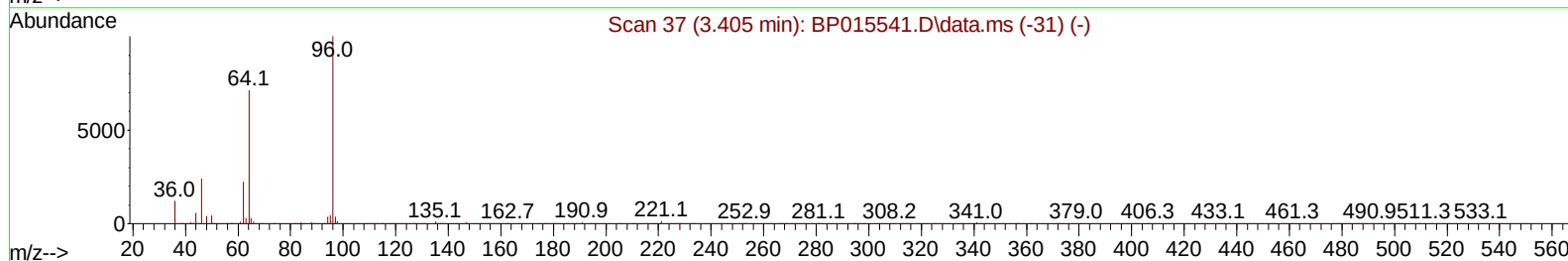
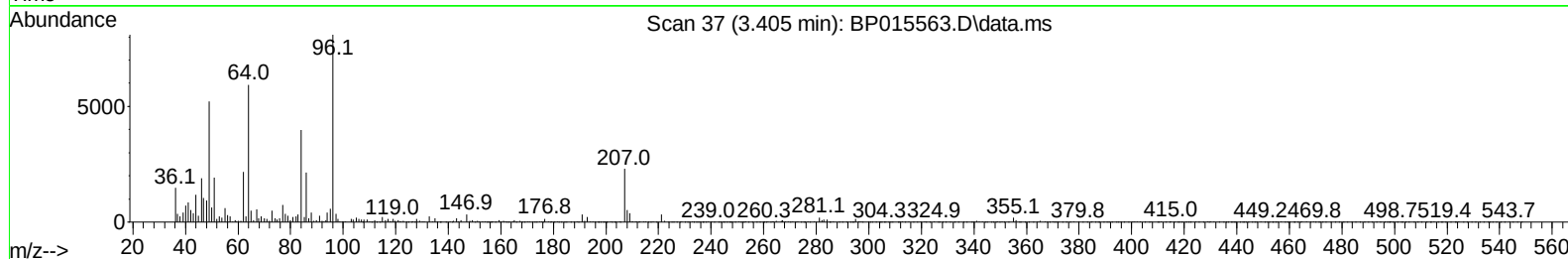
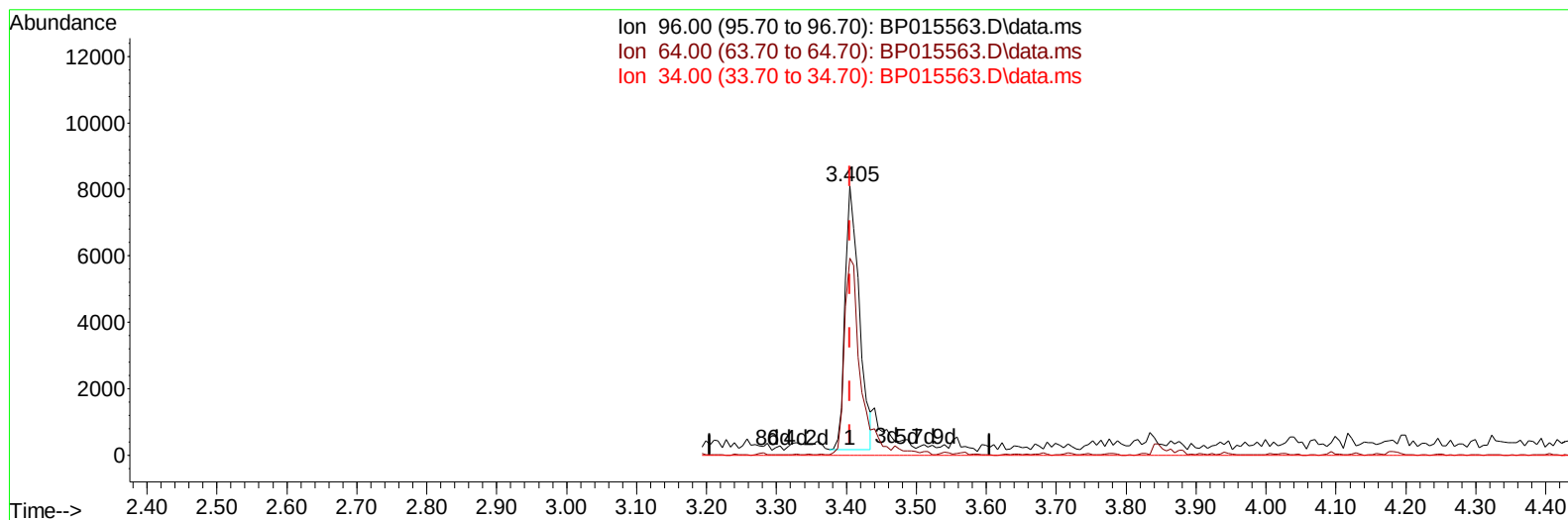
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP061423\
 Data File : BP015563.D
 Acq On : 15 Jun 2023 19:16
 Operator : MA/JU
 Sample : 03062-06
 Misc :
 ALS Vial : 57 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 EZEL2

Manual IntegrationsAPPROVED

Quant Time: Jun 15 22:07:23 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP060723.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Jun 15 05:36:46 2023
 Response via : Initial Calibration

Reviewed By :Yogesh Patel 06/16/2023
 Supervised By :mohammad ahmed 06/16/2023



TIC: BP015563.D\data.ms

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

3.405min (-0.000) 3.86 ng/uL

response 11207

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	75.60	73.25
34.00	0.00	0.00
0.00	0.00	0.00

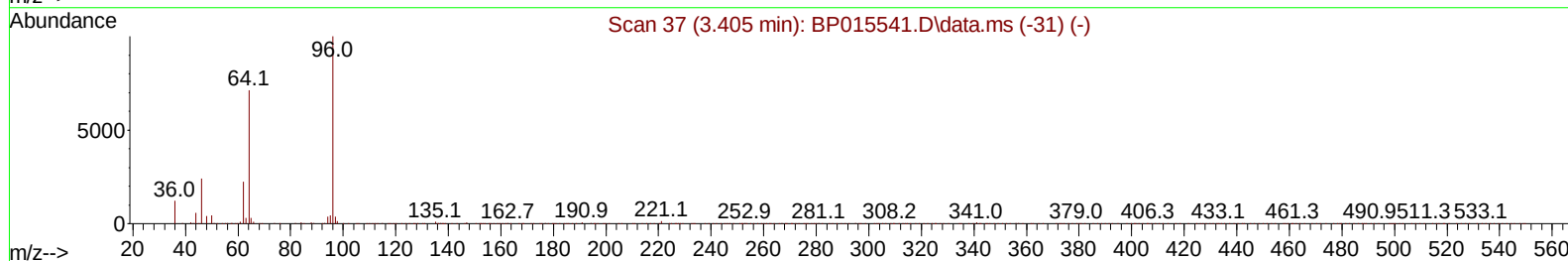
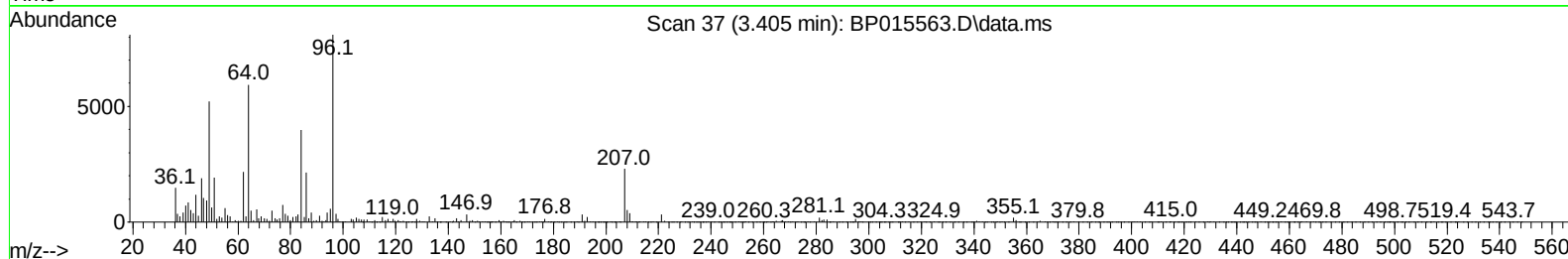
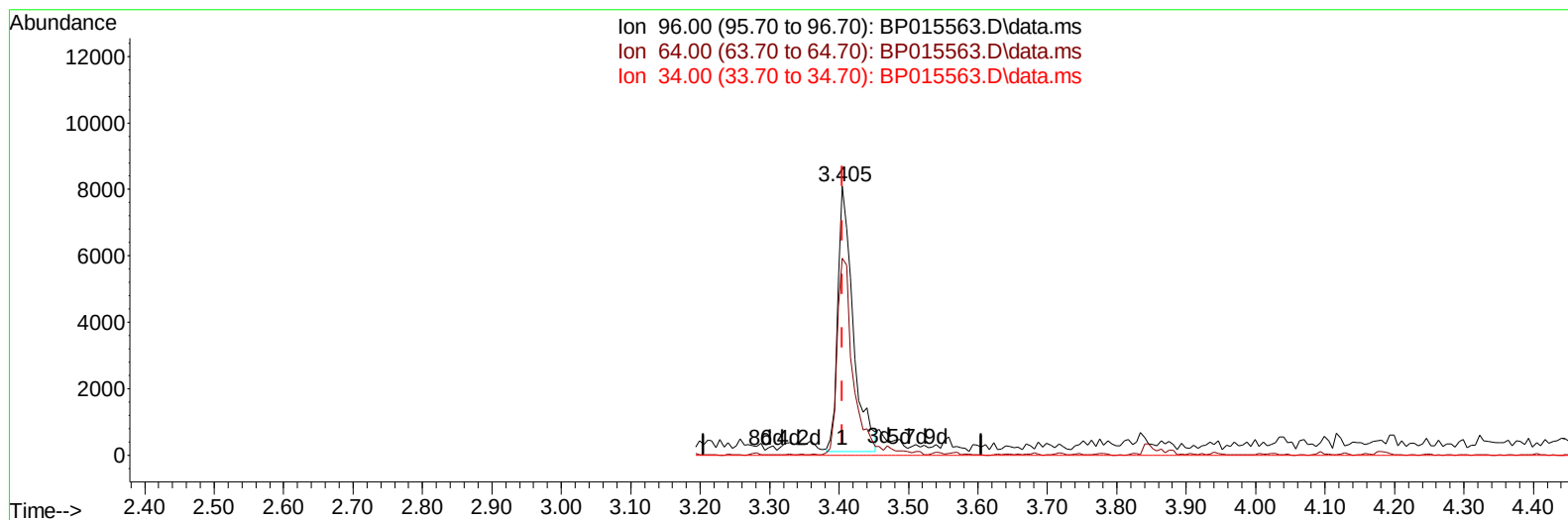
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(3) 1,4-Dioxane-d8 (S)

3.405min (-0.000) 4.23 ng/uL m

response 12282

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	75.60	73.25
34.00	0.00	0.00
0.00	0.00	0.00

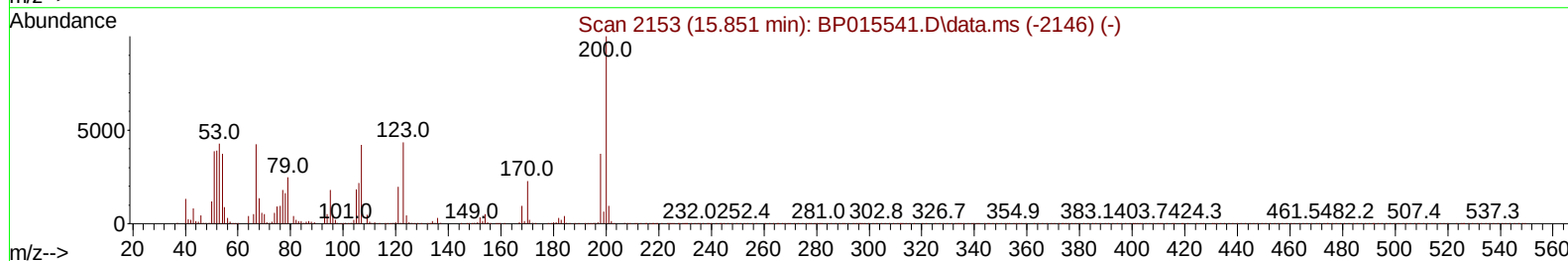
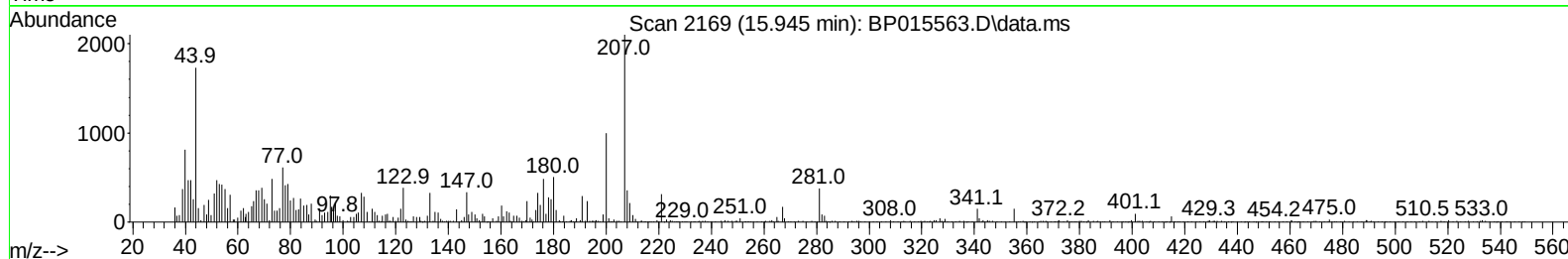
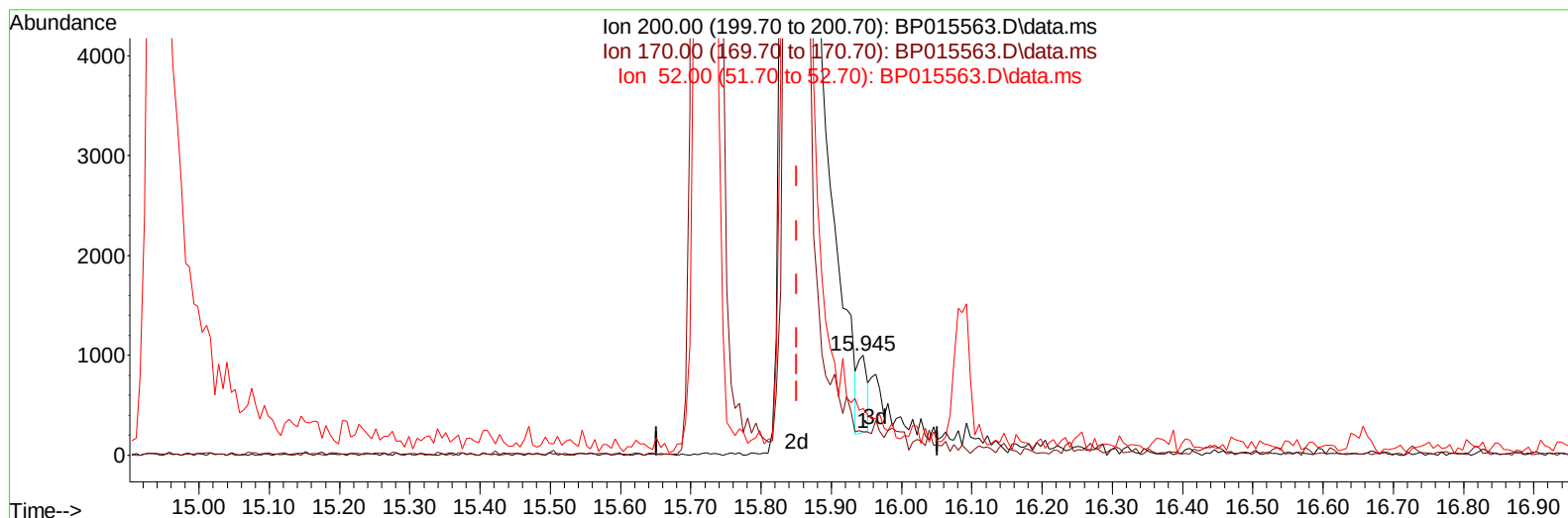
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TIC: BP015563.D\data.ms

(65) 4,6-Dinitro-2-methylphenol-d2 (S)

15.945min (+ 0.094) 0.13 ng/ul

response 709

Ion	Exp%	Act%
200.00	100.00	100.00
170.00	22.30	23.30
52.00	46.20	46.70
0.00	0.00	0.00

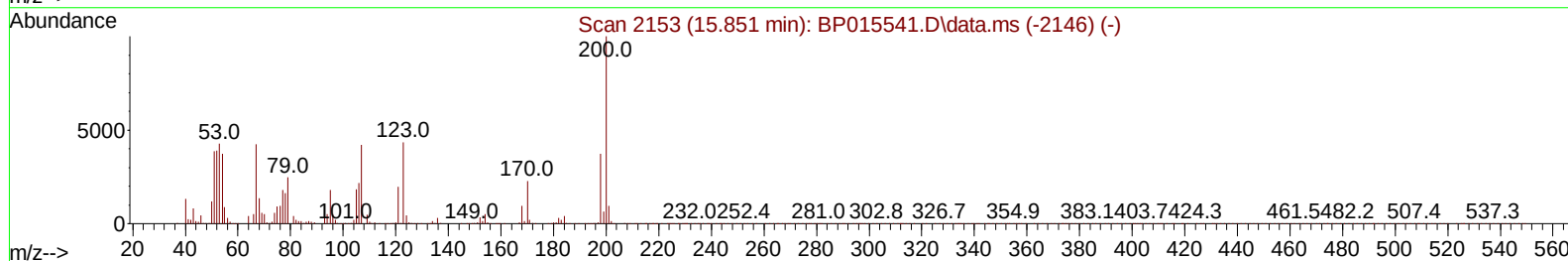
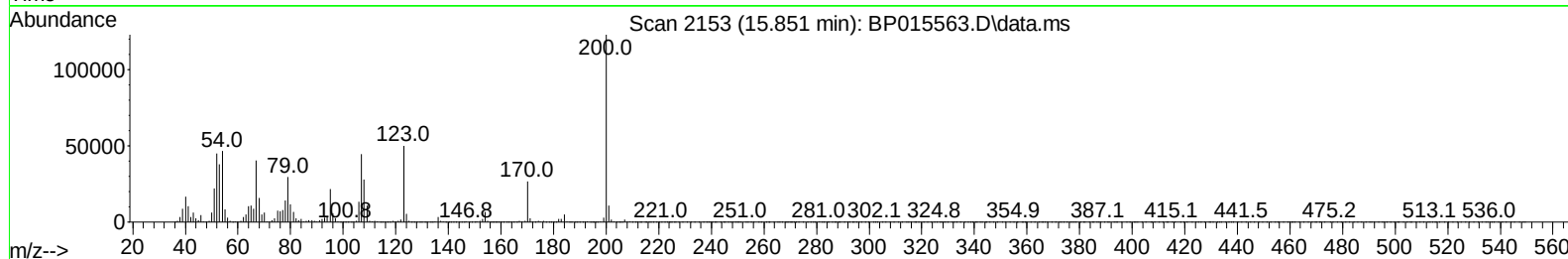
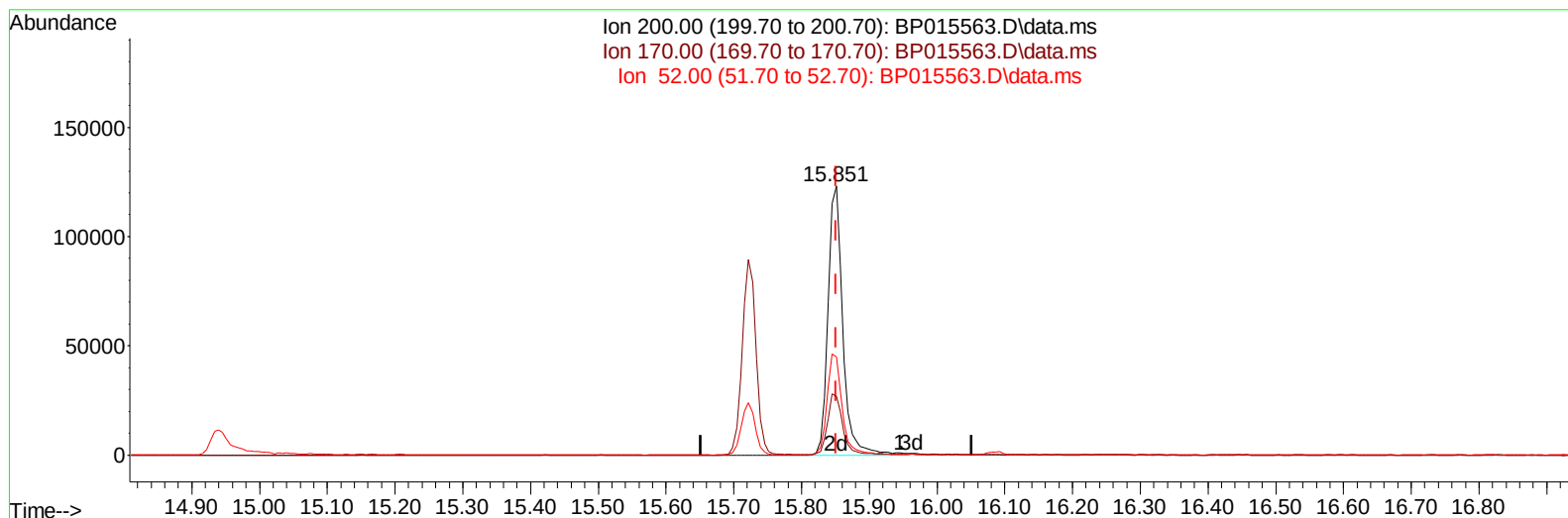
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TIC: BP015563.D\data.ms

(65) 4,6-Dinitro-2-methylphenol-d2 (S)

15.851min (-0.000) 34.25 ng/ul m

response 186032

Ion	Exp%	Act%
200.00	100.00	100.00
170.00	22.30	21.80
52.00	46.20	36.56#
0.00	0.00	0.00

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Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 06/16/2023
 Supervised By :mohammad ahmed 06/16/2023

Quant Time: Jun 15 23:18:48 2023
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.087	152	107405	20.000	ng/ul	0.00
20) Naphthalene-d8	10.910	136	504099	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.728	164	354995	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.486	188	857489	20.000	ng/ul	0.00
79) Chrysene-d12	21.569	240	913291	20.000	ng/ul	0.00
88) Perylene-d12	24.110	264	1006778	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.405	96	12282m	4.234	ng/uL	0.00
4) Pyridine-d5	3.846	84	190212	25.247	ng/ul	0.00
7) Phenol-d5	7.246	99	282364	28.536	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.405	67	172513	29.193	ng/ul	0.00
11) 2-Chlorophenol-d4	7.610	132	207298	28.896	ng/ul	0.00
15) 4-Methylphenol-d8	8.805	113	288819	35.565	ng/ul	0.00
21) Nitrobenzene-d5	9.263	128	134748	34.047	ng/ul	0.00
24) 2-Nitrophenol-d4	9.993	143	163978	36.281	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.534	165	289941	35.168	ng/ul	0.00
31) 4-Chloroaniline-d4	11.057	131	433513	36.632	ng/ul	0.00
46) Dimethylphthalate-d6	14.134	166	1229671	43.927	ng/ul	0.00
49) Acenaphthylene-d8	14.428	160	1314797	42.954	ng/ul	0.00
54) 4-Nitrophenol-d4	14.940	143	177475	35.403	ng/ul	0.00
60) Fluorene-d10	15.722	176	1039292	44.026	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.851	200	186032m	34.246	ng/ul	0.00
73) Anthracene-d10	17.586	188	1743116	44.679	ng/ul	0.00
81) Pyrene-d10	19.810	212	2196303	44.932	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.951	264	2275138	45.038	ng/ul	0.00

Target Compounds Qvalue

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

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