

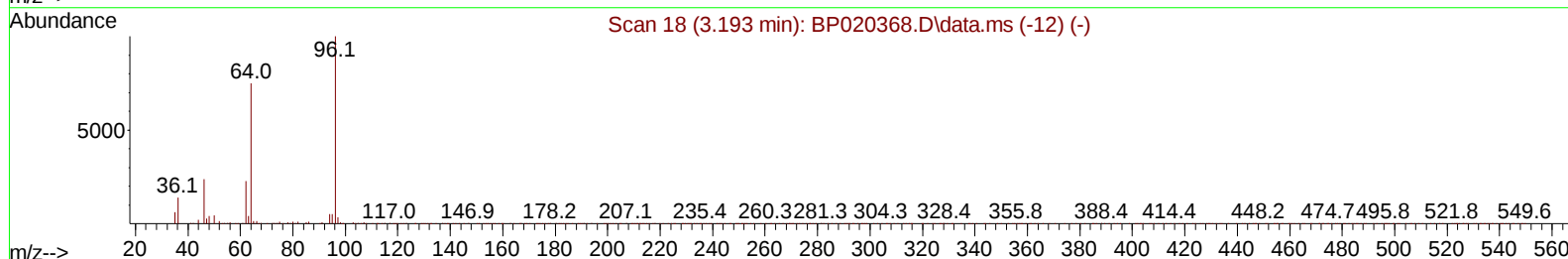
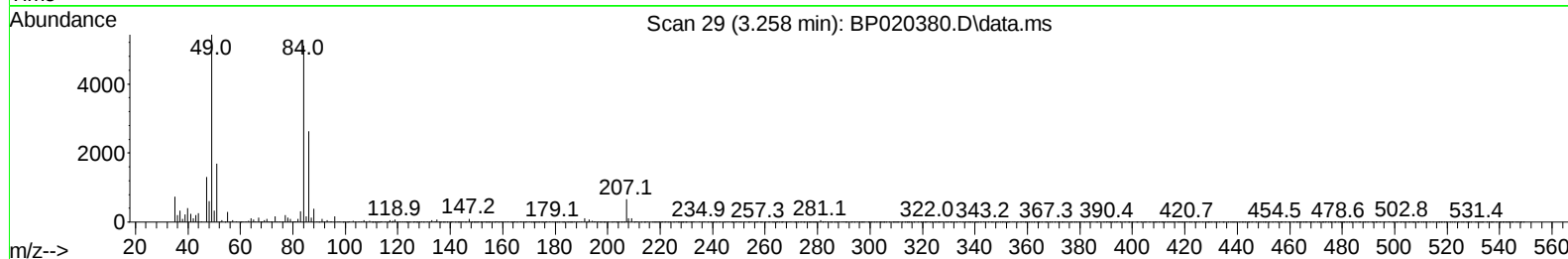
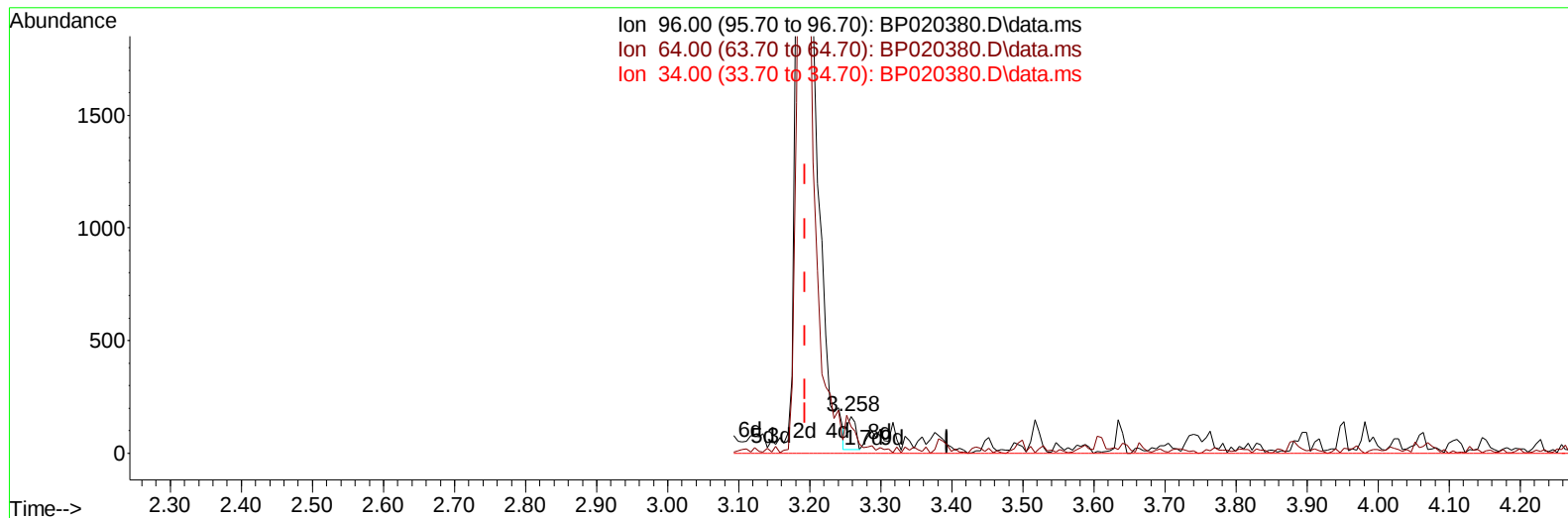
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051424\
 Data File : BP020380.D
 Acq On : 15 May 2024 00:14
 Operator : MA/JU
 Sample : P2372-12
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 A43W2

Manual IntegrationsAPPROVED

Quant Time: May 15 00:41:56 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP050124.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue May 14 23:58:44 2024
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 05/15/2024
 Supervised By :mohammad ahmed 05/17/2024



TIC: BP020380.D\data.ms

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

3.258min (+ 0.065) 0.09 ng/uL

response 129

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	69.80	72.39
34.00	0.00	0.00
0.00	0.00	0.00

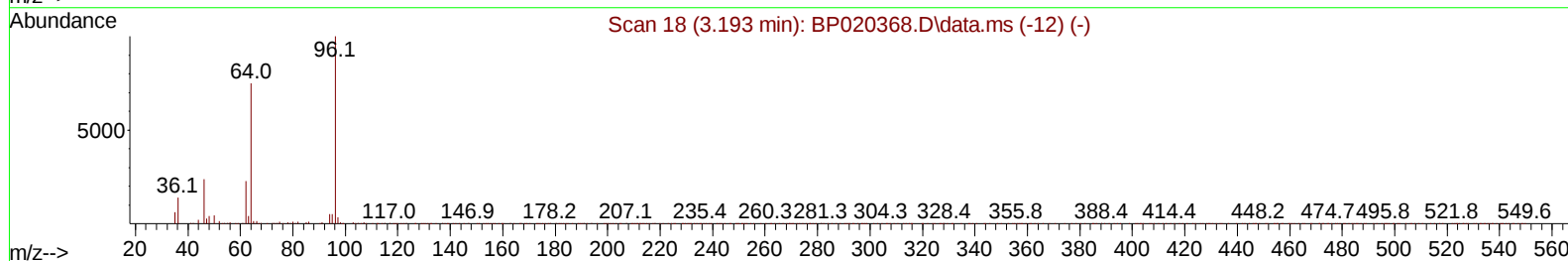
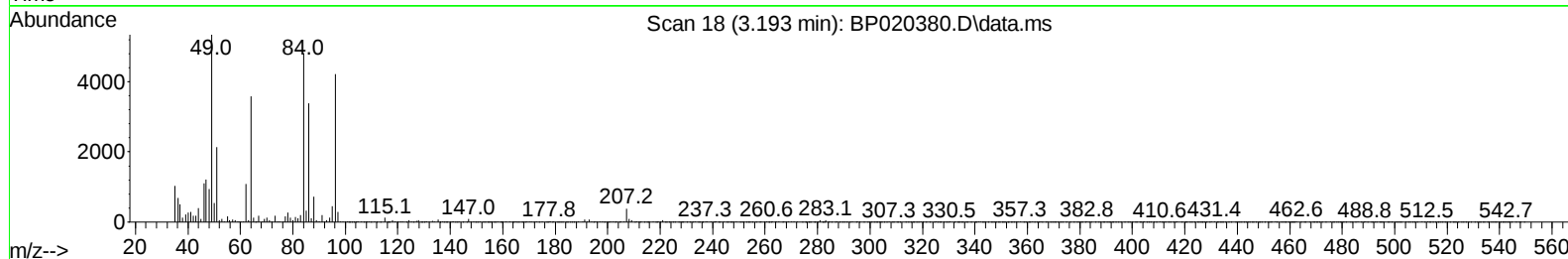
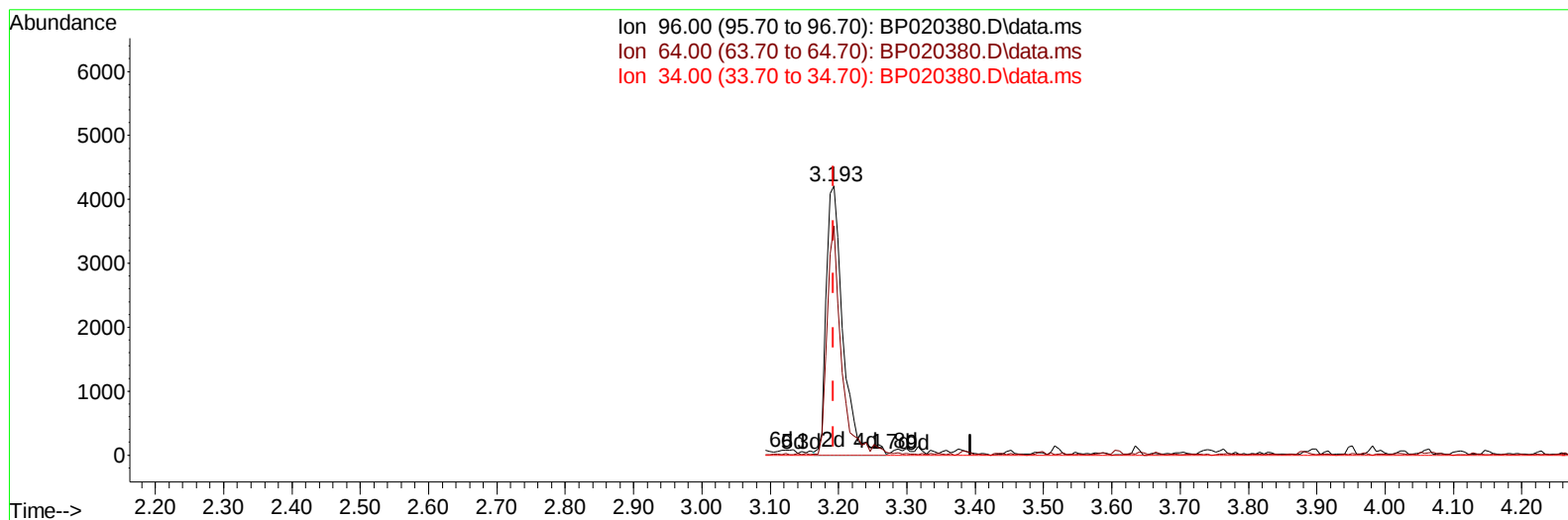
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TIC: BP020380.D\data.ms

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

3.193min (+ 0.000) 4.87 ng/uL m

response 7183

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	69.80	85.31#
34.00	0.00	0.00
0.00	0.00	0.00

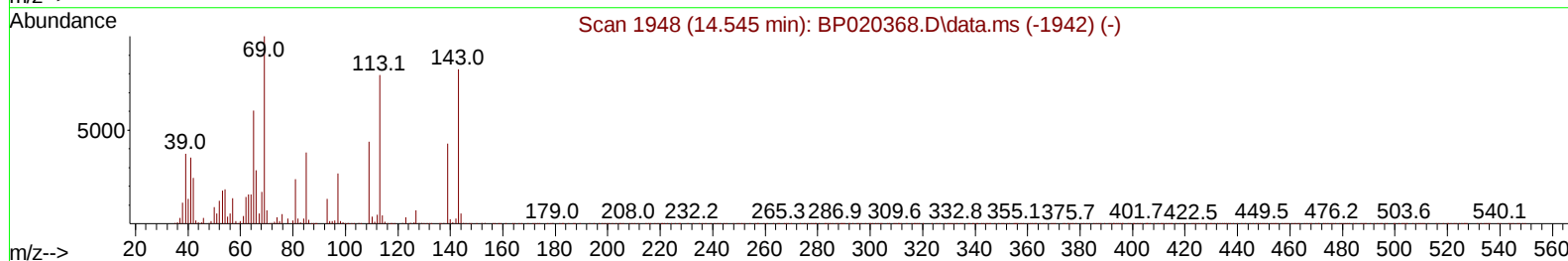
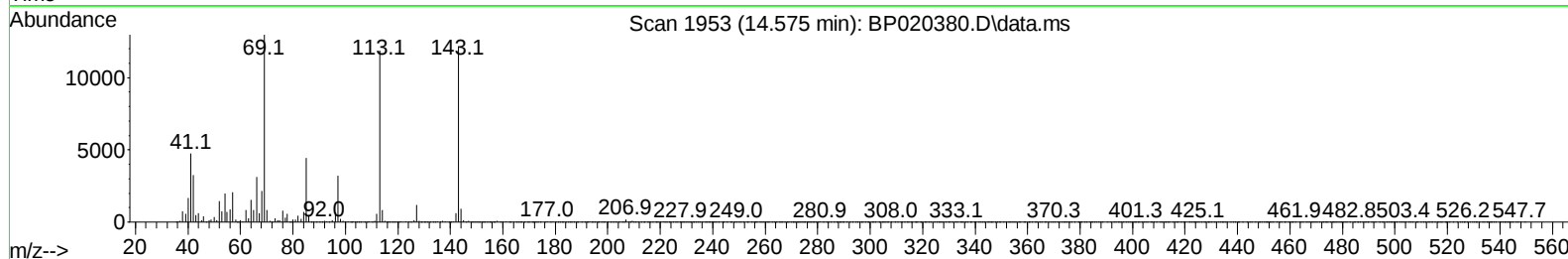
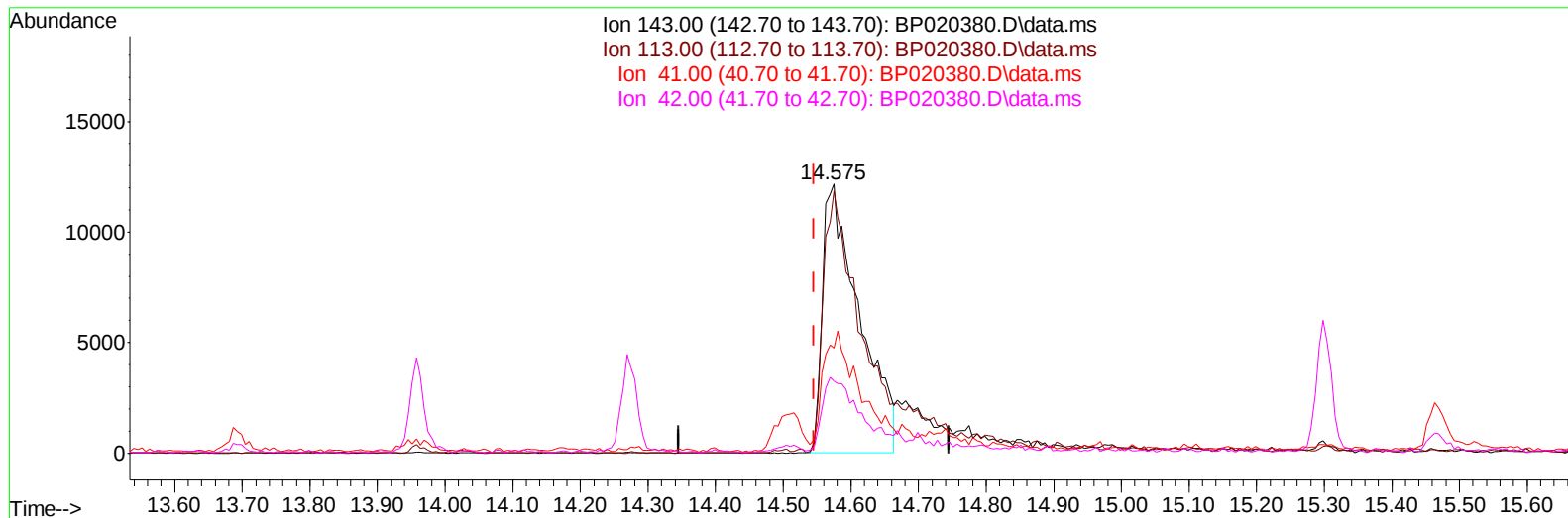
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TIC: BP020380.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.575min (+ 0.030) 20.95 ng/u1

response 46063

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	101.40	97.61
41.00	50.30	39.13#
42.00	32.50	26.72

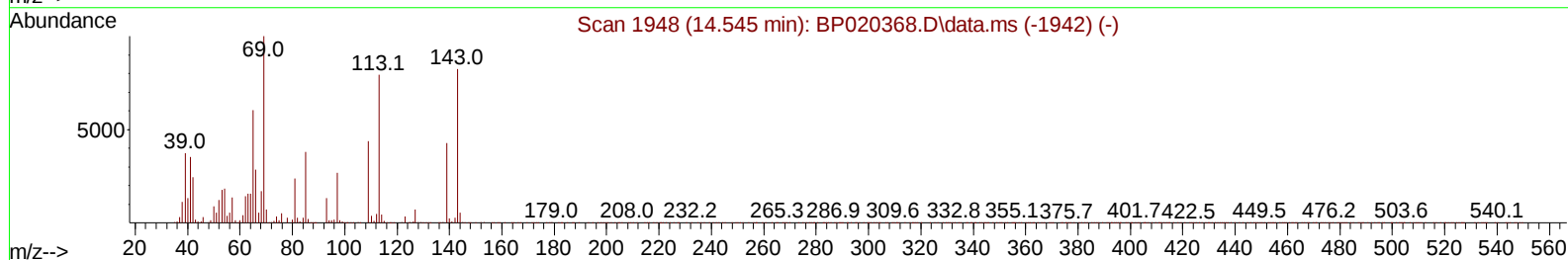
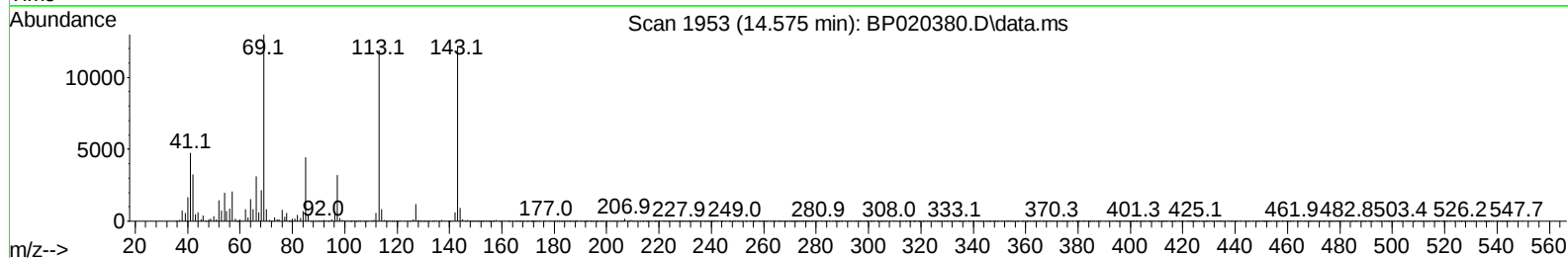
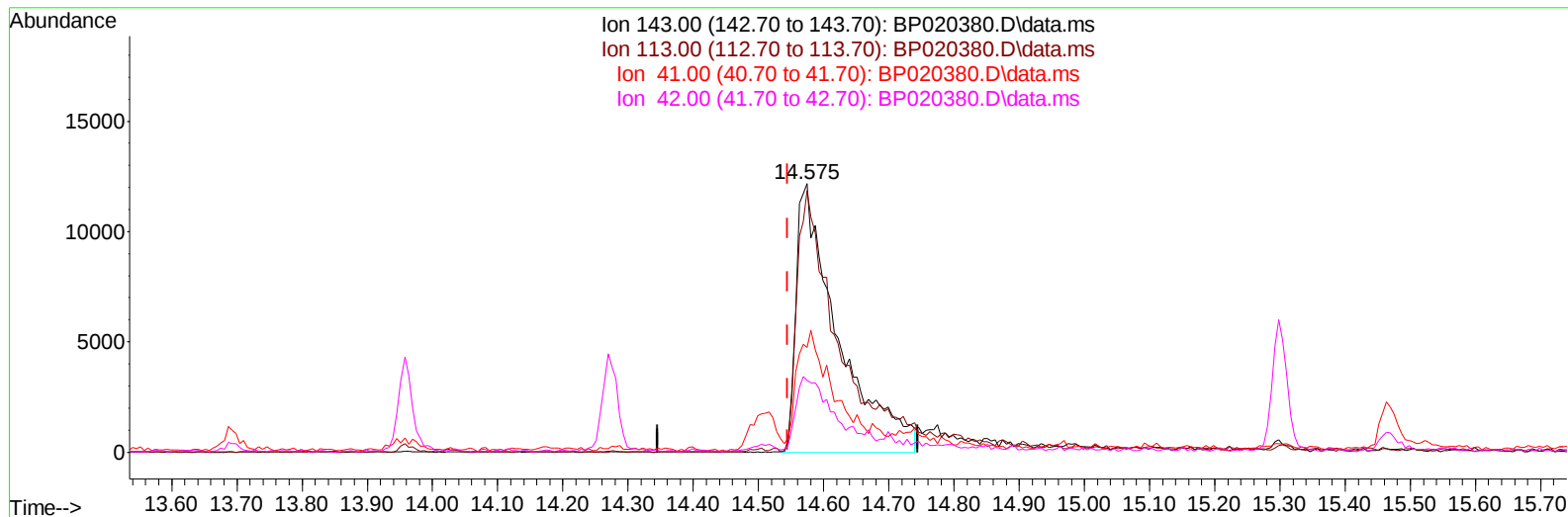
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ALS Vial : 14 Sample Multiplier: 1

Instrument :
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TIC: BP020380.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.575min (+ 0.030) 24.57 ng/ul m

response 54022

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	101.40	97.61
41.00	50.30	39.13#
42.00	32.50	26.72

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Instrument :
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Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 05/15/2024
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.593	152	62354	20.000	ng/ul	0.00
20) Naphthalene-d8	10.363	136	266549	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.275	164	195886	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.110	188	483505	20.000	ng/ul	0.00
79) Chrysene-d12	21.598	240	483542	20.000	ng/ul	0.00
88) Perylene-d12	24.939	264	554232	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.193	96	7183m	4.865	ng/uL	0.00
4) Pyridine-d5	3.599	84	91180	21.254	ng/ul	0.00
7) Phenol-d5	6.787	99	130602	24.389	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.952	67	78283	24.065	ng/ul	0.00
11) 2-Chlorophenol-d4	7.134	132	102181	26.692	ng/ul	0.00
15) 4-Methylphenol-d8	8.305	113	103186	23.834	ng/ul	0.00
21) Nitrobenzene-d5	8.758	128	50139	25.120	ng/ul	0.00
24) 2-Nitrophenol-d4	9.469	143	57922	25.341	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.010	165	99737	23.911	ng/ul	0.00
31) 4-Chloroaniline-d4	10.534	131	122615	21.004	ng/ul	0.00
46) Dimethylphthalate-d6	13.693	166	422849	29.560	ng/ul	0.00
49) Acenaphthylene-d8	13.957	160	430100	27.218	ng/ul	0.00
54) 4-Nitrophenol-d4	14.575	143	54022m	24.572	ng/ul	0.03
60) Fluorene-d10	15.298	176	352679	29.559	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.469	200	56794	18.511	ng/ul	0.01
73) Anthracene-d10	17.216	188	599852	28.678	ng/ul	0.00
81) Pyrene-d10	19.610	212	752467	26.621	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.710	264	762235	28.444	ng/ul	0.00

Target Compounds Qvalue

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

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