

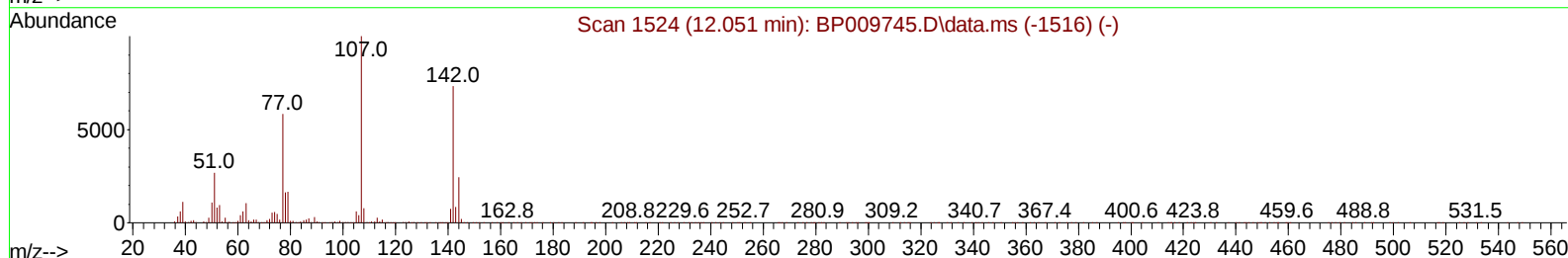
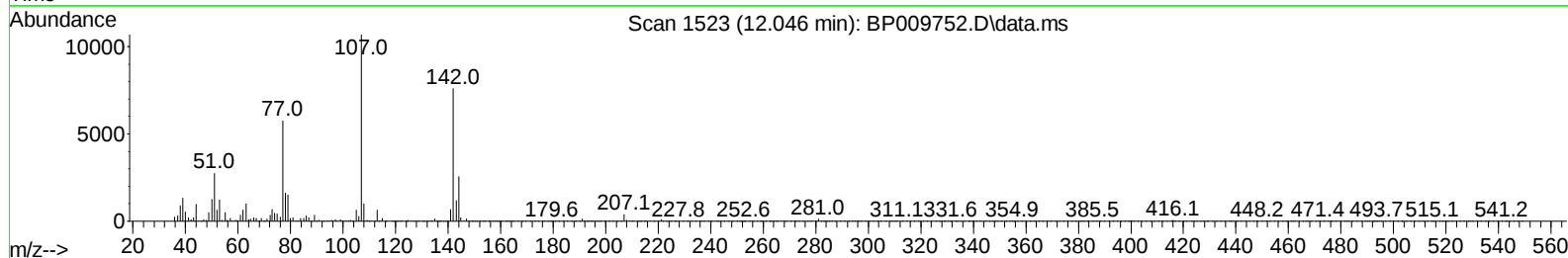
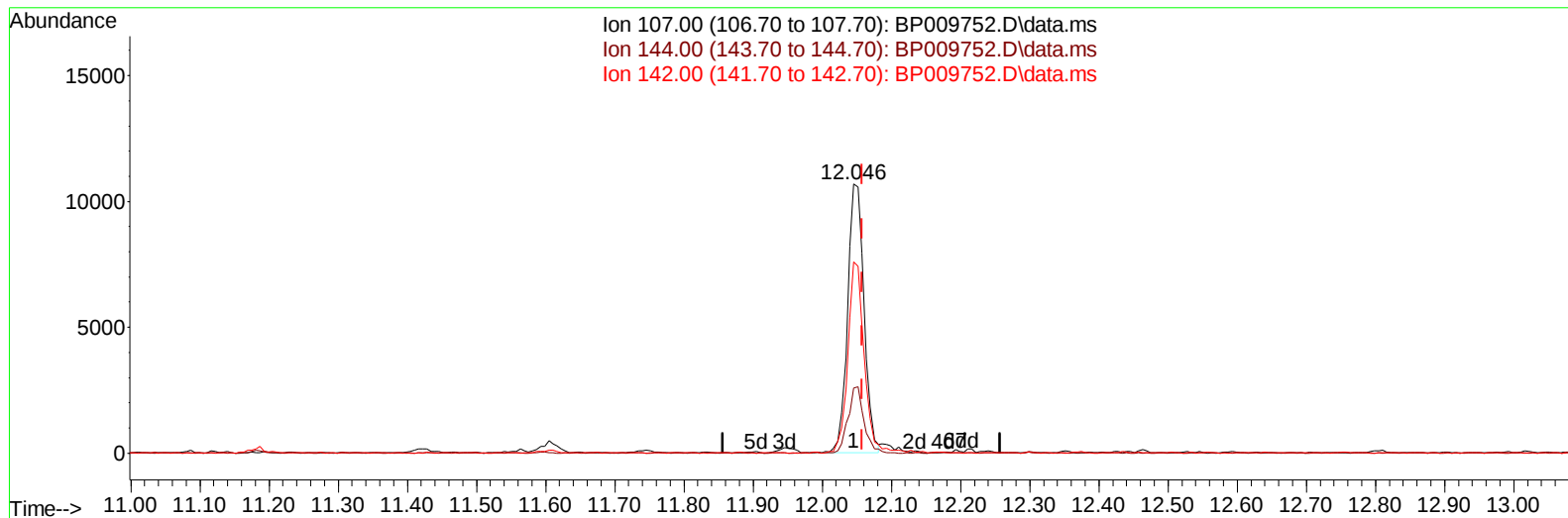
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041122\
 Data File : BP009752.D
 Acq On : 11 Apr 2022 15:23
 Operator : CG/JU
 Sample : N2338-09
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0HY0

Manual IntegrationsAPPROVED

Quant Time: Apr 12 01:35:33 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP040722.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Apr 07 19:16:52 2022
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 04/12/2022
 Supervised By :Yogesh Patel 04/13/2022



TIC: BP009752.D\data.ms

(35) 4-Chloro-3-methylphenol (C)

12.046min (-0.012) 1.89 ng/u1

response 17586

Ion	Exp%	Act%
107.00	100.00	100.00
144.00	19.80	24.07#
142.00	48.60	71.18#
0.00	0.00	0.00

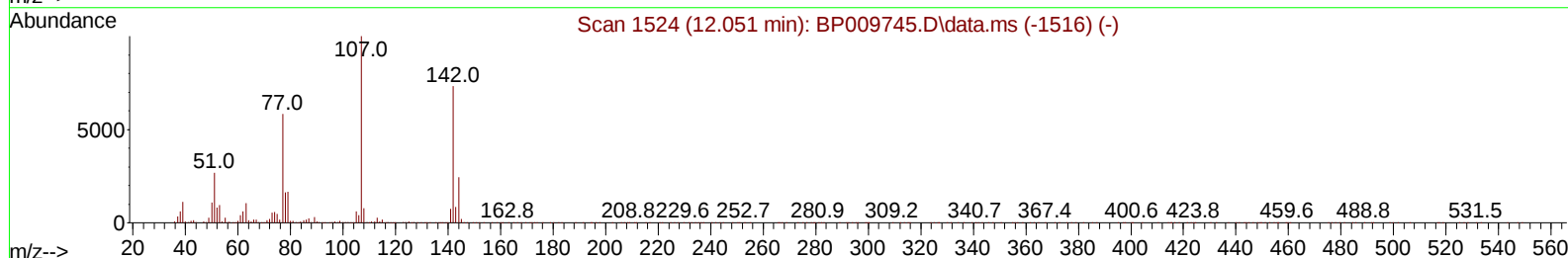
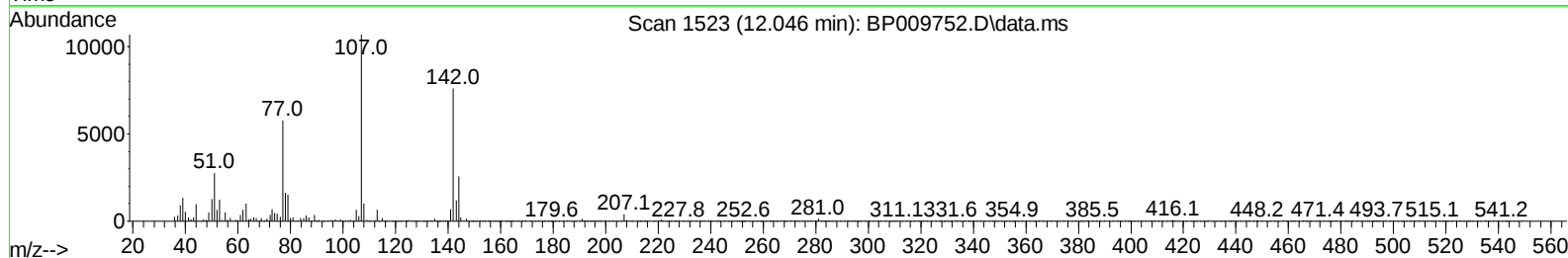
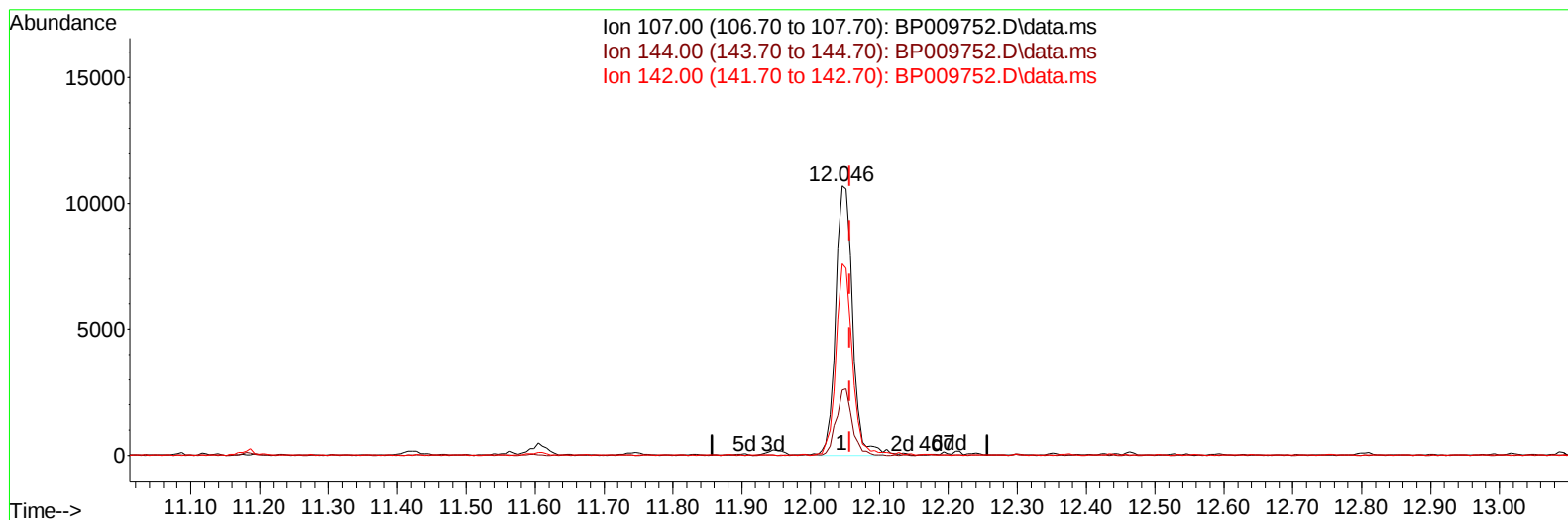
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TIC: BP009752.D\data.ms

(35) 4-Chloro-3-methylphenol (C)

12.046min (-0.012) 1.94 ng/ul m

response 18019

Ion	Exp%	Act%
107.00	100.00	100.00
144.00	19.80	24.07#
142.00	48.60	71.18#
0.00	0.00	0.00

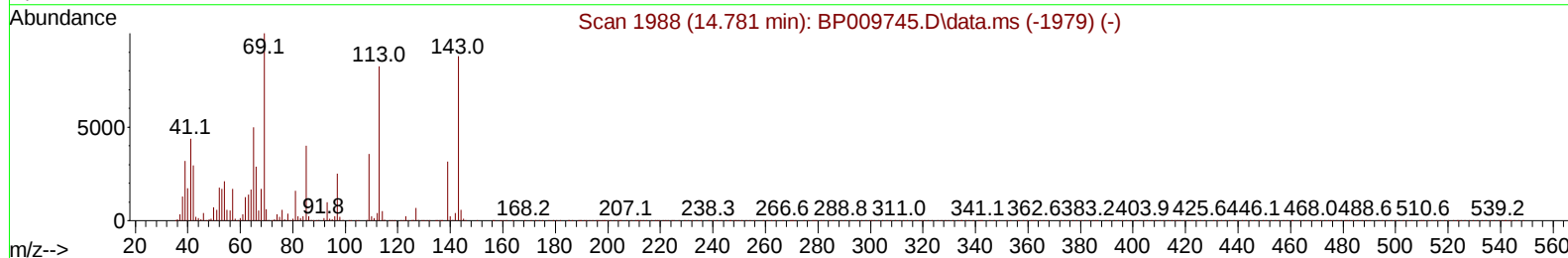
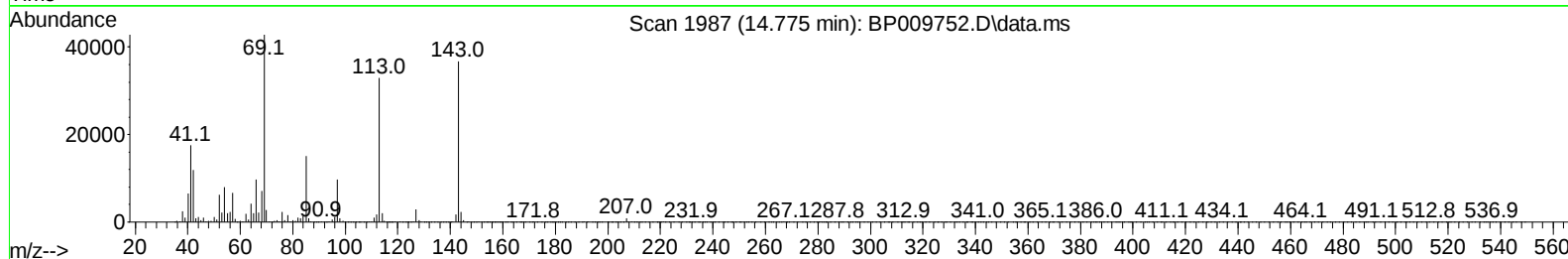
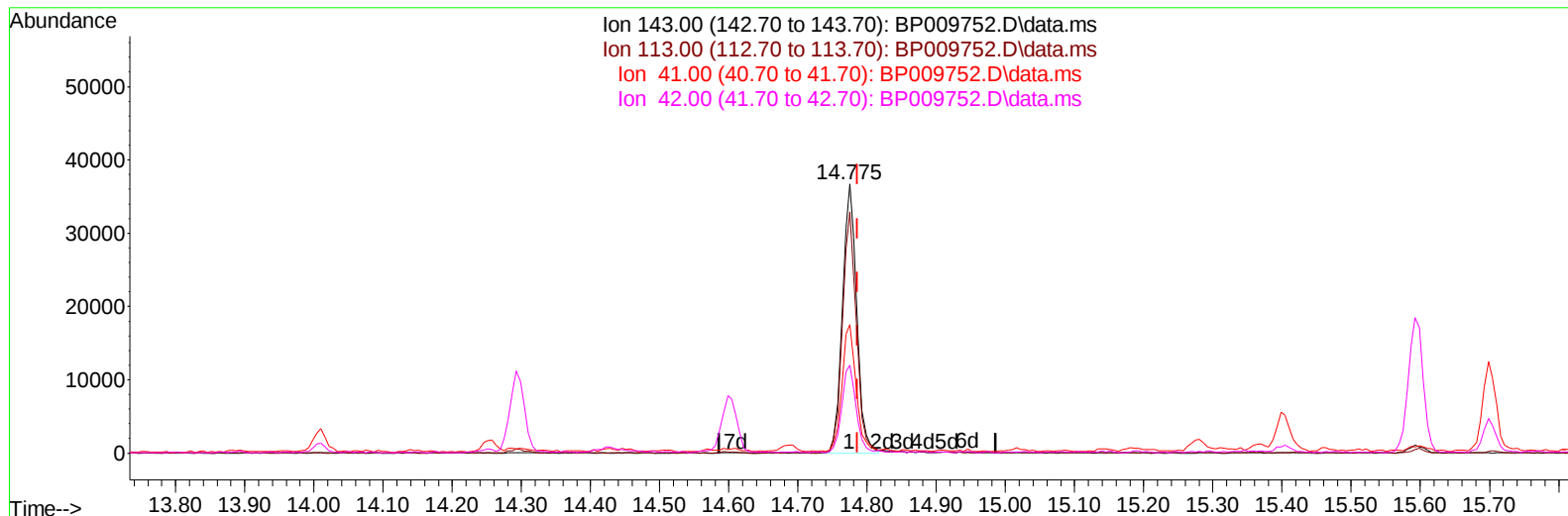
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 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
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TIC: BP009752.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.775min (-0.012) 11.00 ng/ul

response 52730

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	132.70	89.64#
41.00	23.20	47.70#
42.00	18.00	32.59#

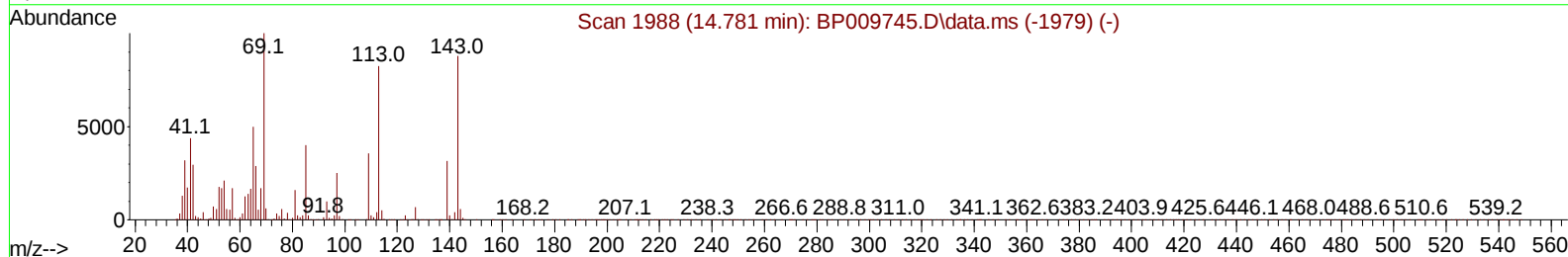
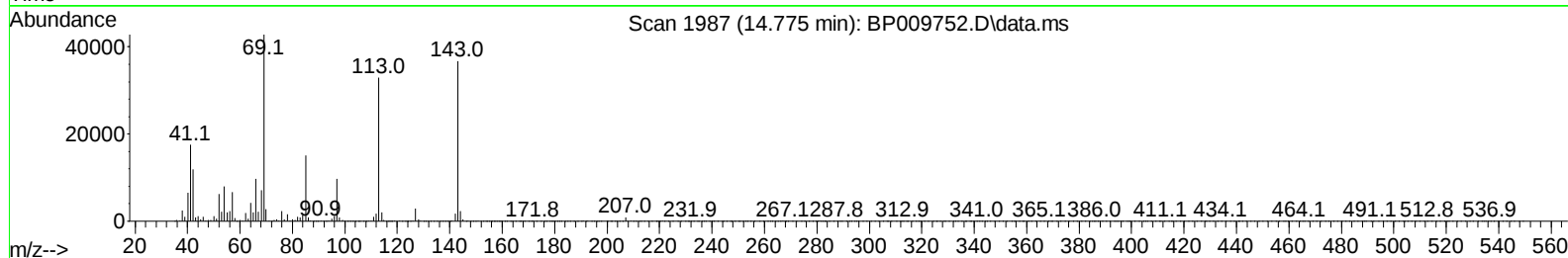
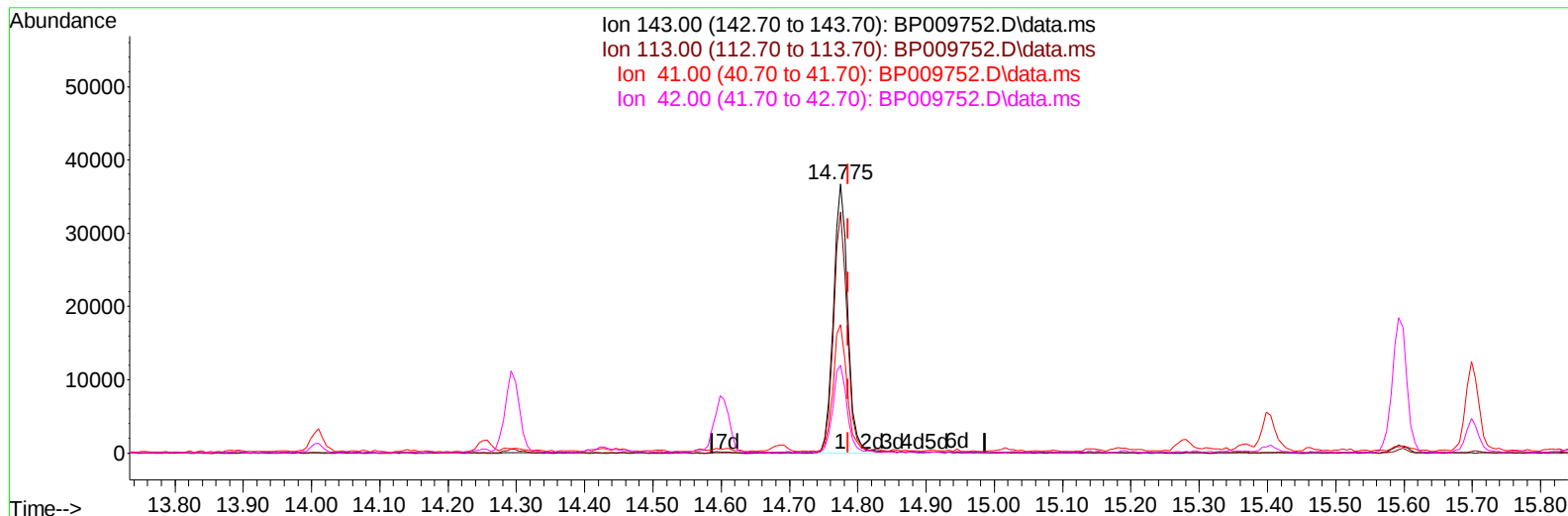
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 Sample : N2338-09
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_P
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TIC: BP009752.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.775min (-0.012) 11.09 ng/ul m

response 53139

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	132.70	89.64#
41.00	23.20	47.70#
42.00	18.00	32.59#

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 Misc :
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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	7.969	152	112210	20.000	ng/ul	-0.01
20) Naphthalene-d8	10.781	136	478238	20.000	ng/ul	-0.01
38) Acenaphthene-d10	14.604	164	336422	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.351	188	751440	20.000	ng/ul	#-0.01
79) Chrysene-d12	21.433	240	841663	20.000	ng/ul	#-0.01
88) Perylene-d12	23.904	264	835159	20.000	ng/ul	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.387	96	13558	5.392	ng/uL	0.00
4) Pyridine-d5	3.805	84	88488	12.765	ng/ul	0.00
7) Phenol-d5	7.122	99	89692	9.105	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.293	67	226597	38.637	ng/ul	-0.01
11) 2-Chlorophenol-d4	7.499	132	224543	30.217	ng/ul	0.00
15) 4-Methylphenol-d8	8.669	113	164562	20.011	ng/ul	-0.01
21) Nitrobenzene-d5	9.128	128	145329	42.140	ng/ul	-0.01
24) 2-Nitrophenol-d4	9.852	143	151705	40.123	ng/ul	-0.01
28) 2,4-Dichlorophenol-d3	10.393	165	268079	33.883	ng/ul	-0.01
31) 4-Chloroaniline-d4	10.910	131	368005	32.657	ng/ul	0.00
46) Dimethylphthalate-d6	14.010	166	1088755	41.200	ng/ul	-0.01
49) Acenaphthylene-d8	14.298	160	1201600	38.380	ng/ul	0.00
54) 4-Nitrophenol-d4	14.775	143	53139m	11.088	ng/ul	-0.01
60) Fluorene-d10	15.592	176	924081	41.568	ng/ul	-0.01
65) 4,6-Dinitro-2-methylph...	15.698	200	183969	41.327	ng/ul	-0.01
73) Anthracene-d10	17.451	188	1653931	45.970	ng/ul	-0.01
81) Pyrene-d10	19.680	212	2048383	44.561	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.751	264	2069088	47.574	ng/ul	-0.01
Target Compounds						
35) 4-Chloro-3-methylphenol	12.046	107	18019m	1.939	ng/ul	Qvalue

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

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