

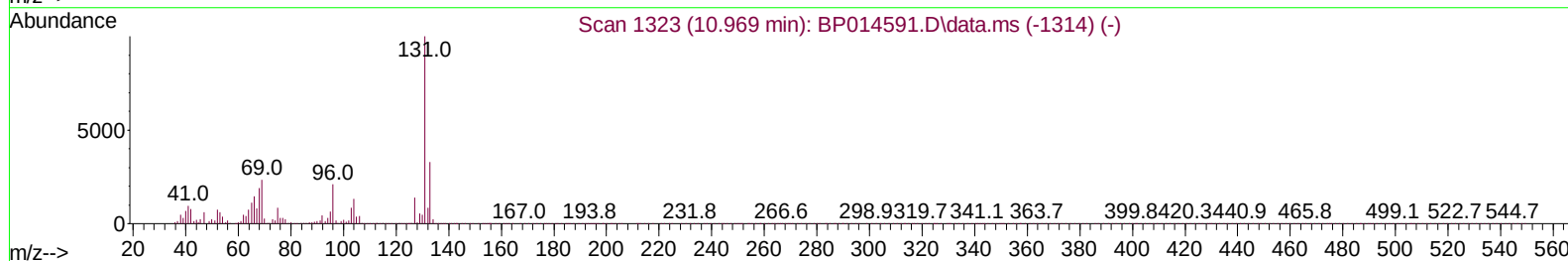
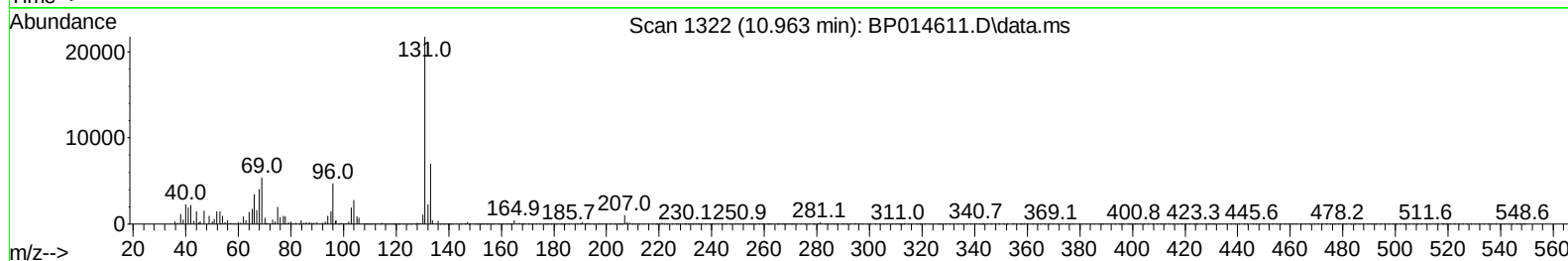
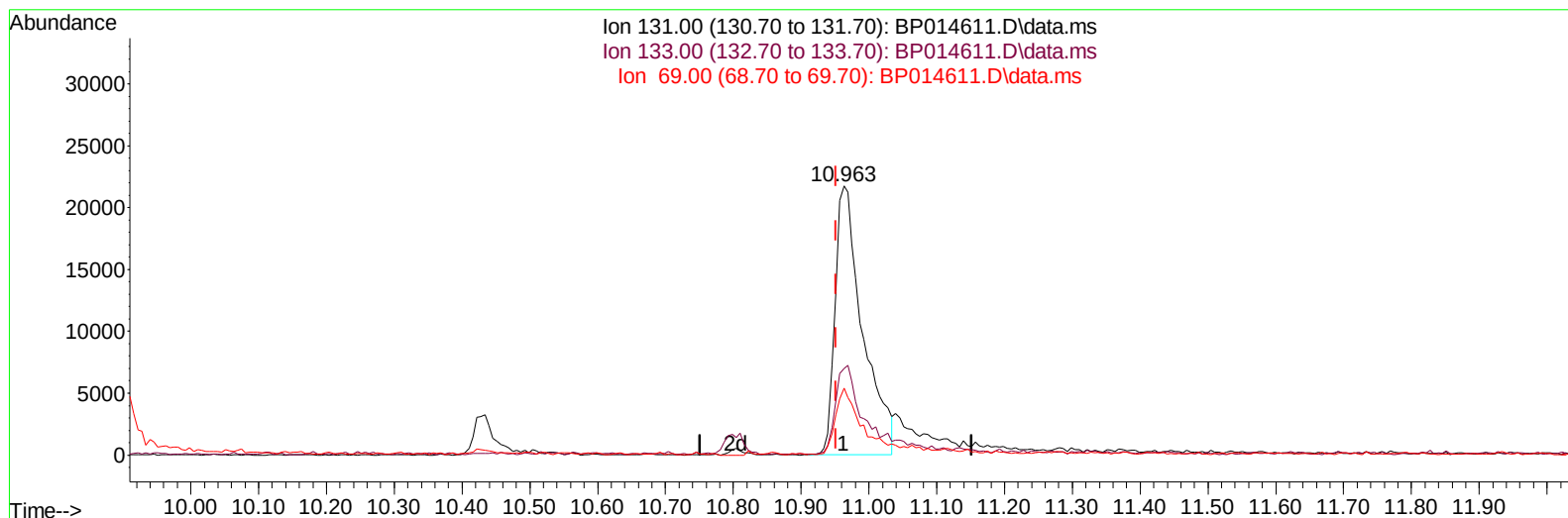
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP040723\
Data File : BP014611.D
Acq On : 07 Apr 2023 18:59
Operator : CG/JU
Sample : 02225-15
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0AB7

Manual IntegrationsAPPROVED

Reviewed By :Christian Giraldo 04/10/2023
Supervised By :Jagrut Upadhyay 04/10/2023

Quant Time: Apr 08 01:50:12 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP032823.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri Apr 07 00:45:34 2023
Response via : Initial Calibration



TIC: BP014611.D\data.ms

(31) 4-Chloroaniline-d4 (S)

10.963min (+ 0.011) 6.76 ng/ul

response 61455

Ion	Exp%	Act%
131.00	100.00	100.00
133.00	32.20	32.24
69.00	21.30	24.88
0.00	0.00	0.00

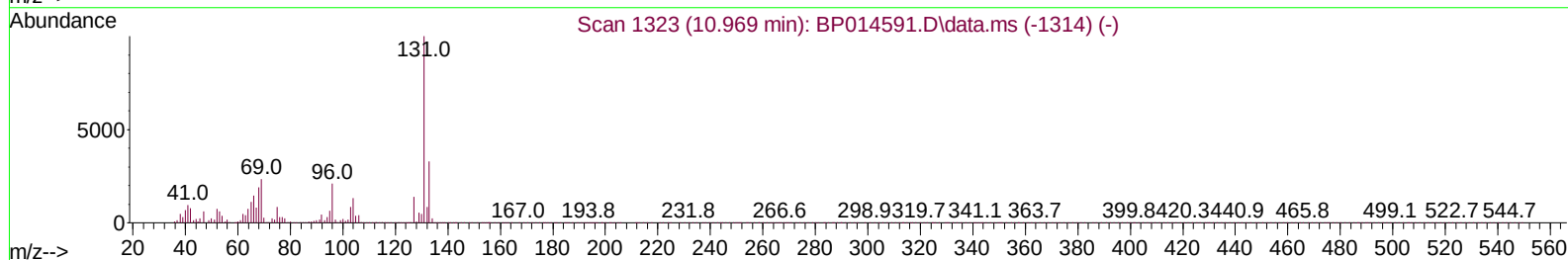
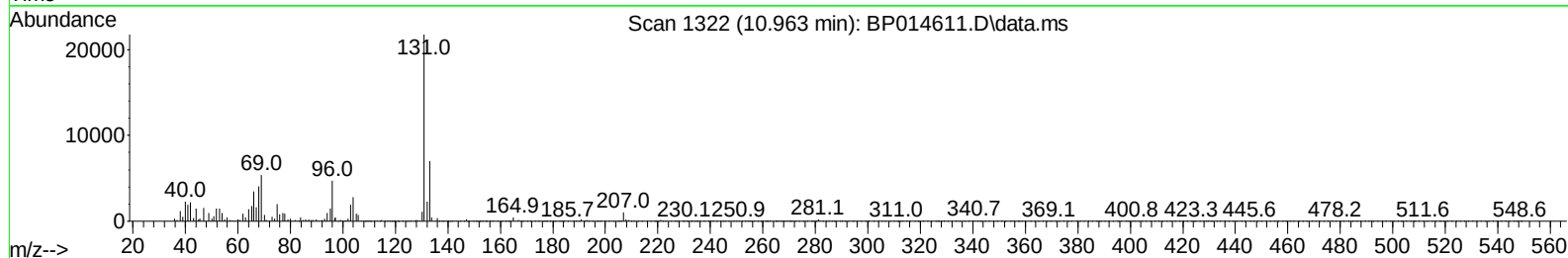
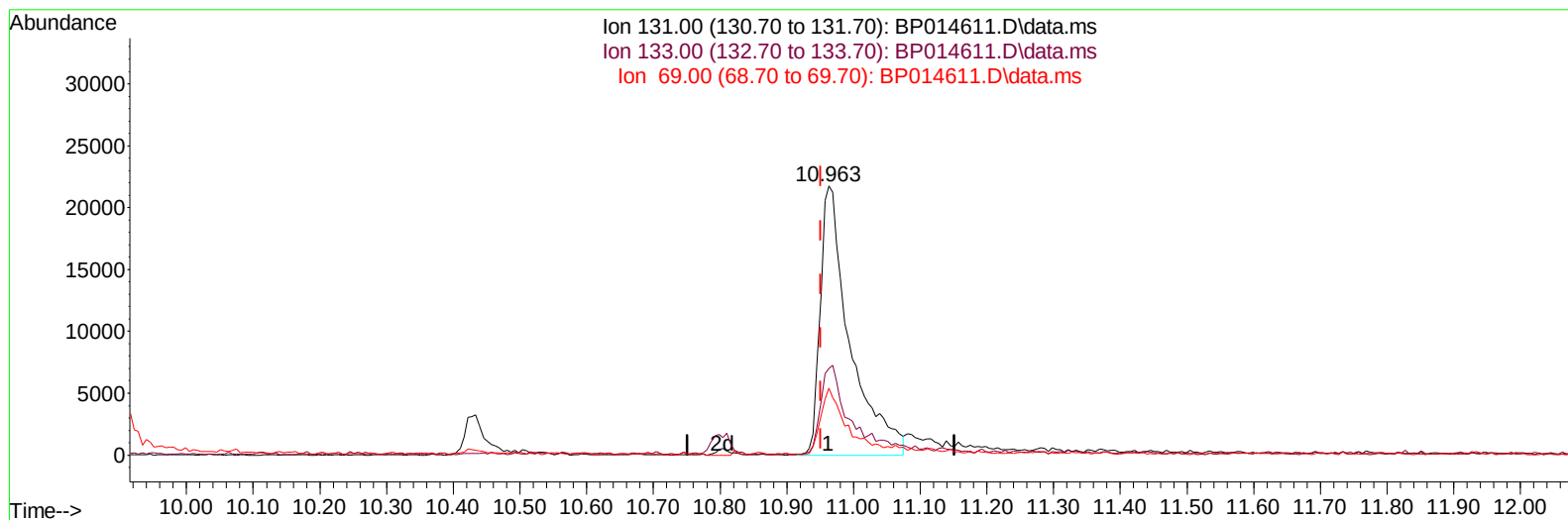
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TIC: BP014611.D\data.ms

(31) 4-Chloroaniline-d4 (S)

10.963min (+ 0.011) 7.40 ng/ul m

response 67206

Ion	Exp%	Act%
131.00	100.00	100.00
133.00	32.20	32.24
69.00	21.30	24.88
0.00	0.00	0.00

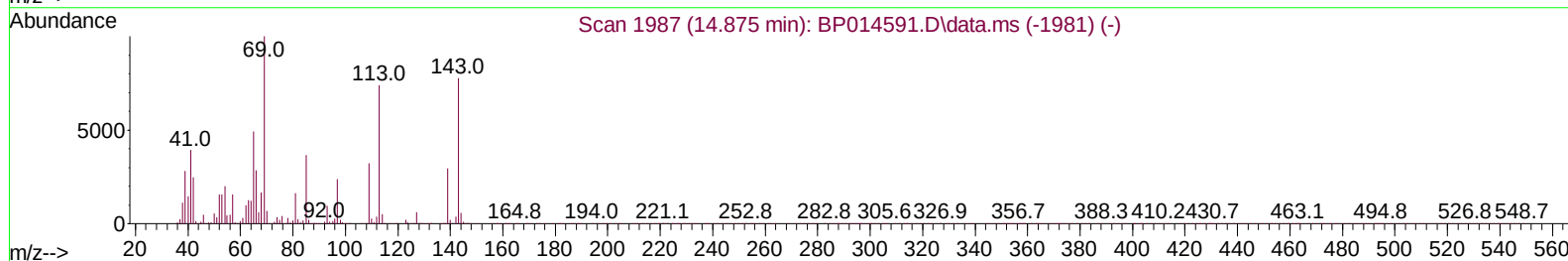
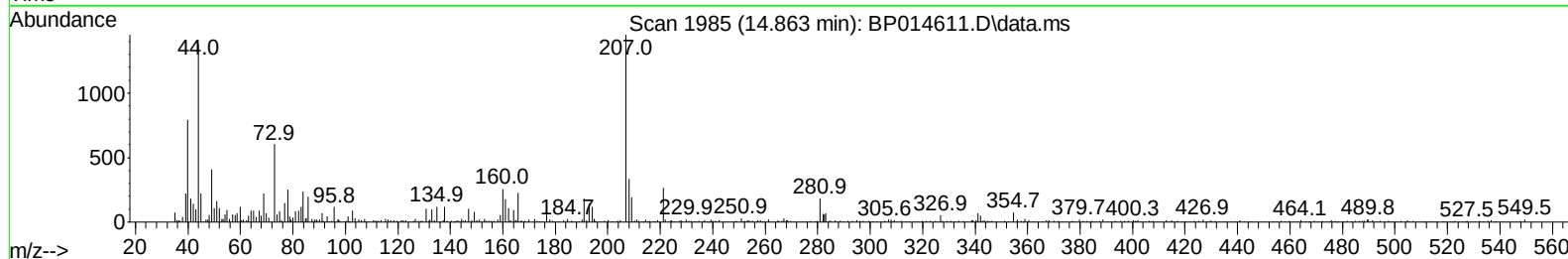
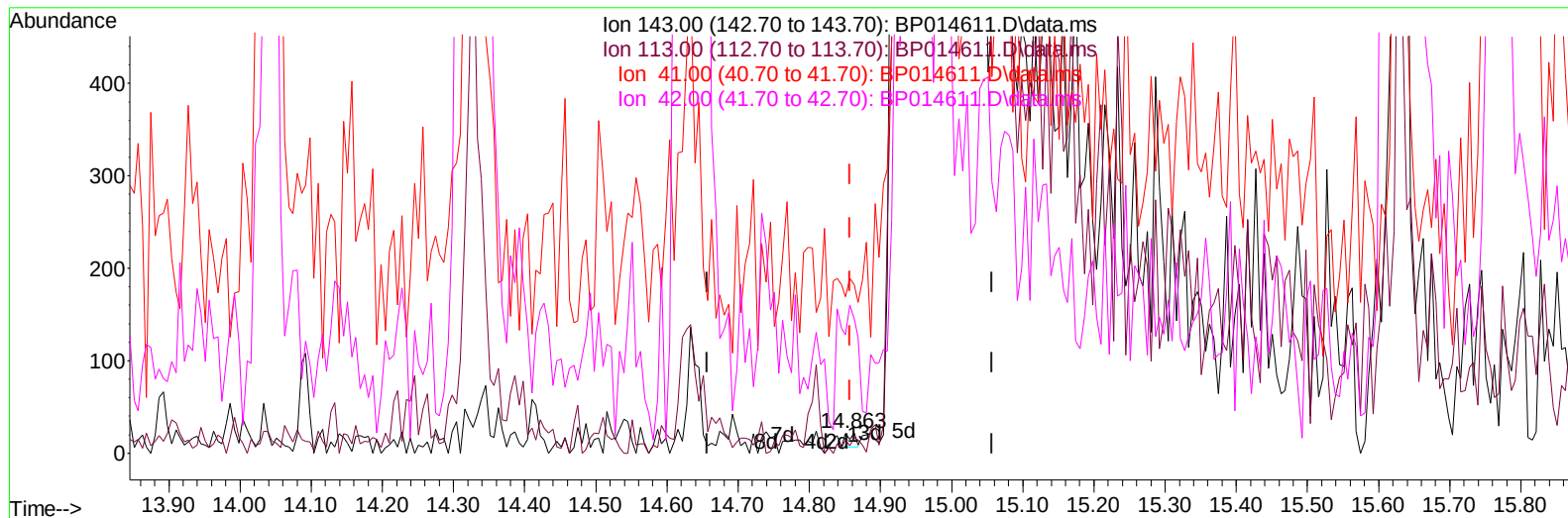
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TIC: BP014611.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.863min (+ 0.006) 0.00 ng/ul

response 14

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	96.80	77.27#
41.00	44.20	827.27#
42.00	31.50	659.09#

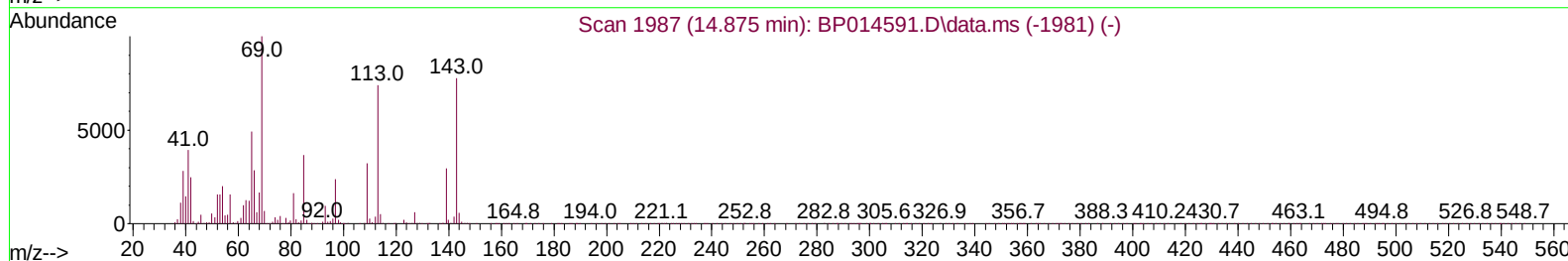
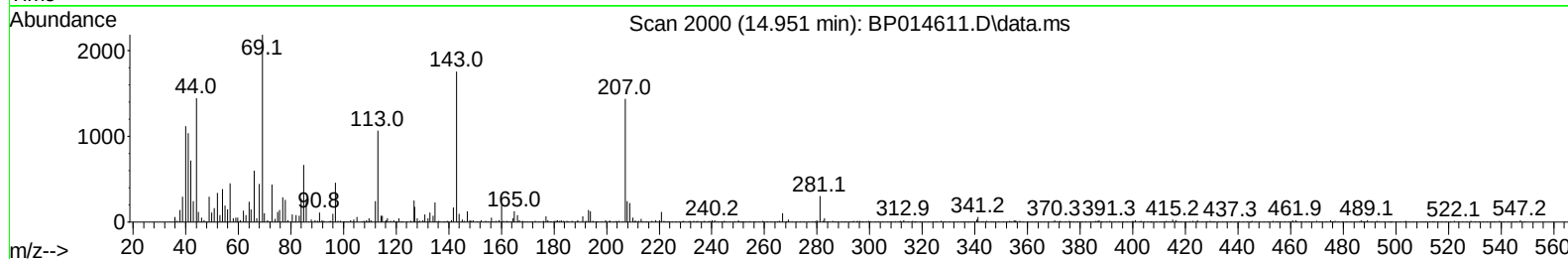
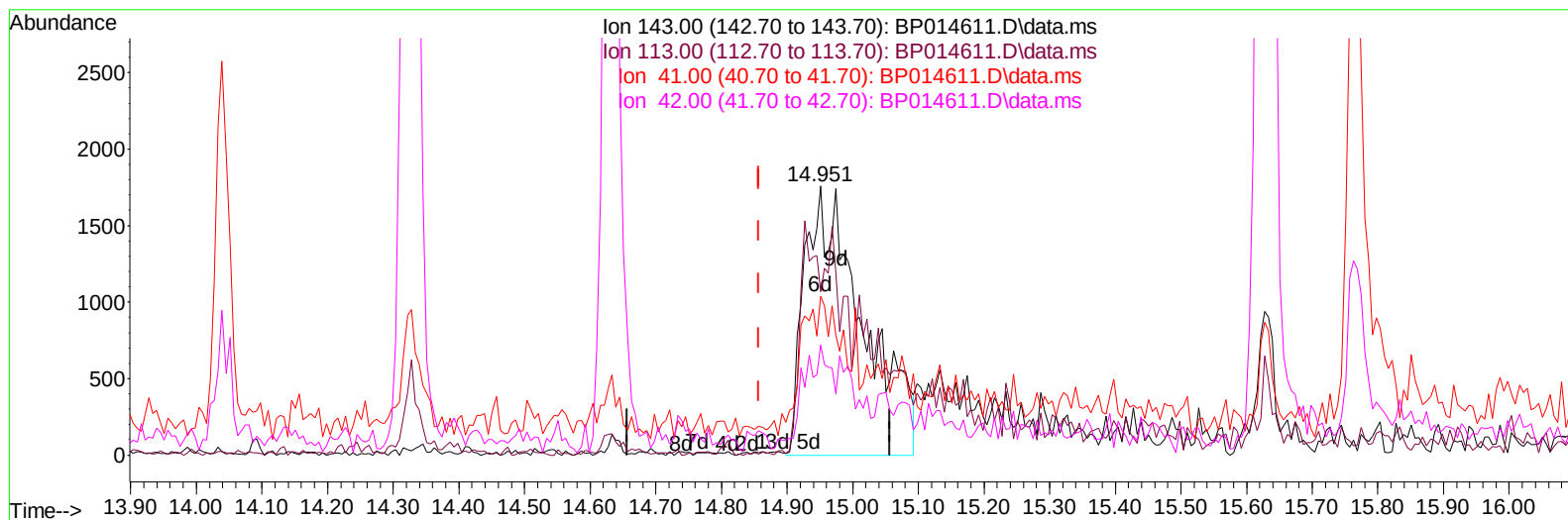
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TIC: BP014611.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.951min (+ 0.094) 3.29 ng/ul m

response 10737

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	96.80	60.75#
41.00	44.20	59.10#
42.00	31.50	41.01#

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 Misc :
 ALS Vial : 4 Sample Multiplier: 1

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.981	152	97679	20.000	ng/ul	0.00
20) Naphthalene-d8	10.798	136	409807	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.633	164	266223	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.398	188	591251	20.000	ng/ul	0.00
79) Chrysene-d12	21.486	240	622613	20.000	ng/ul	0.00
88) Perylene-d12	23.986	264	720683	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.363	96	13344	5.298	ng/uL	0.00
4) Pyridine-d5	0.000	84	0d	0.000	ng/ul	
7) Phenol-d5	7.157	99	53267	5.913	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.304	67	223280	38.631	ng/ul	0.00
11) 2-Chlorophenol-d4	7.510	132	182257	27.992	ng/ul	0.00
15) 4-Methylphenol-d8	8.704	113	111789	15.384	ng/ul	0.00
21) Nitrobenzene-d5	9.157	128	117492	35.514	ng/ul	0.00
24) 2-Nitrophenol-d4	9.881	143	125888	33.231	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.428	165	189058	27.403	ng/ul	0.00
31) 4-Chloroaniline-d4	10.963	131	67206m	7.395	ng/ul	0.01
46) Dimethylphthalate-d6	14.039	166	809484	37.501	ng/ul	0.00
49) Acenaphthylene-d8	14.327	160	884286	36.919	ng/ul	0.00
54) 4-Nitrophenol-d4	14.951	143	10737m	3.293	ng/ul	0.09
60) Fluorene-d10	15.627	176	685741	38.095	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.763	200	67588	15.536	ng/ul	0.00
73) Anthracene-d10	17.498	188	1109093	38.761	ng/ul	0.00
81) Pyrene-d10	19.727	212	1389503	33.225	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.827	264	1506244	39.338	ng/ul	0.00
Target Compounds						
2) 1,4-Dioxane	3.399	88	3808	1.376	ng/uL#	Qvalue 91

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

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