

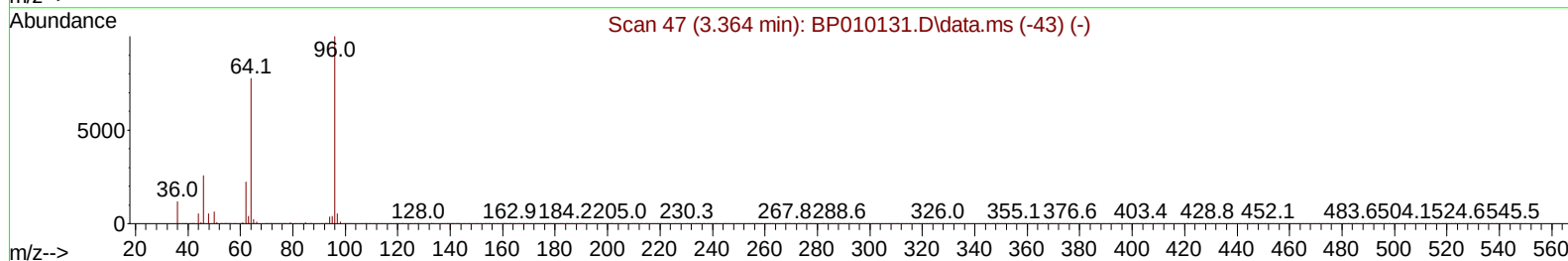
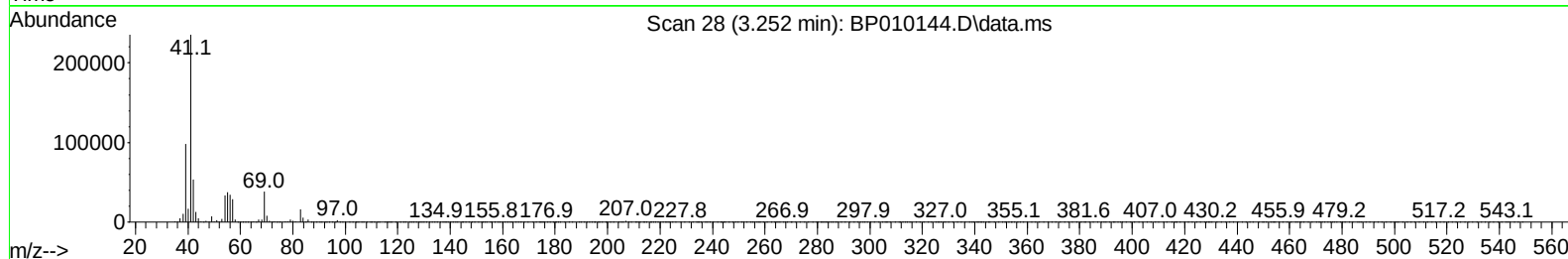
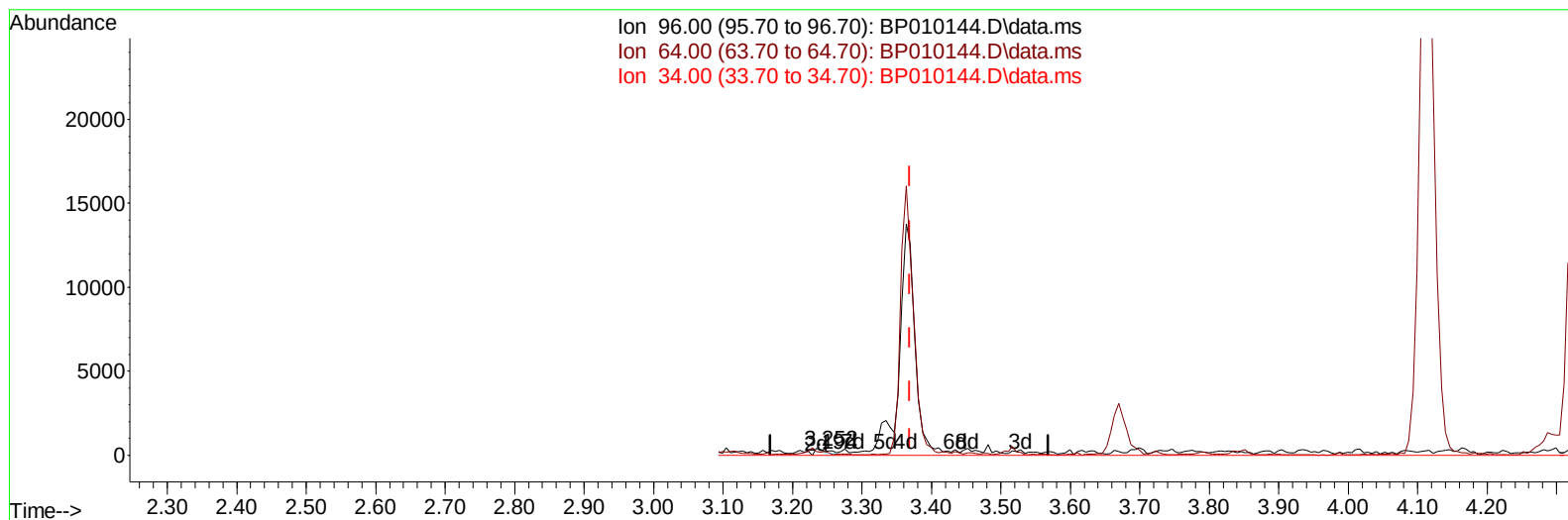
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042922\
 Data File : BP010144.D
 Acq On : 29 Apr 2022 11:03
 Operator : CG/JU
 Sample : N2587-02
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 EXET0

Manual IntegrationsAPPROVED

Quant Time: Apr 29 11:33:32 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP042122.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Apr 21 14:49:38 2022
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 05/02/2022
 Supervised By :Yogesh Patel 05/05/2022



TIC: BP010144.D\data.ms

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

3.252min (-0.117) 0.04 ng/uL

response	141	
Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	50.96
34.00	0.00	0.00
0.00	0.00	0.00

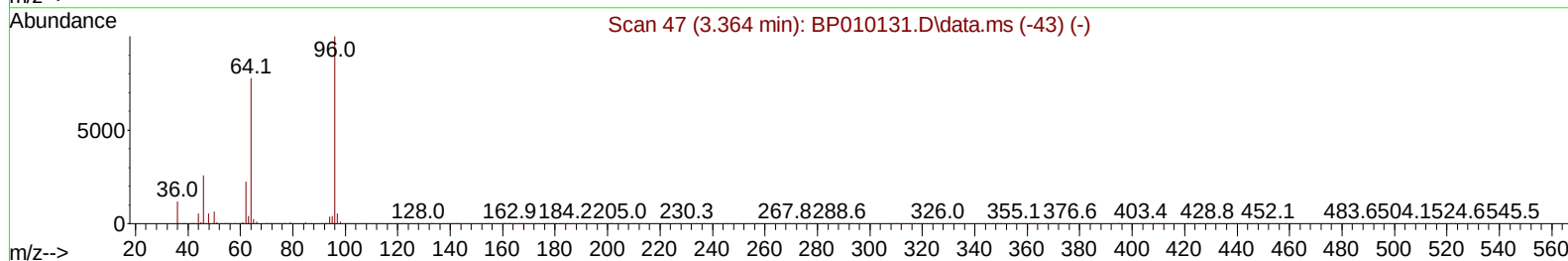
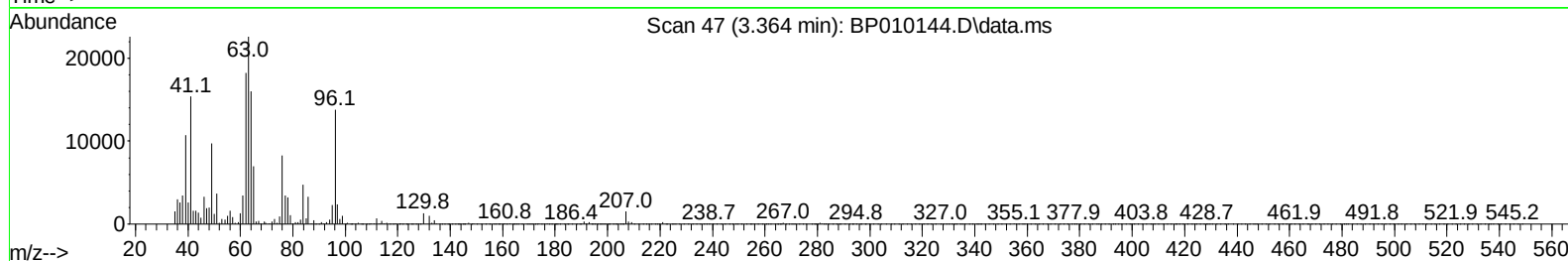
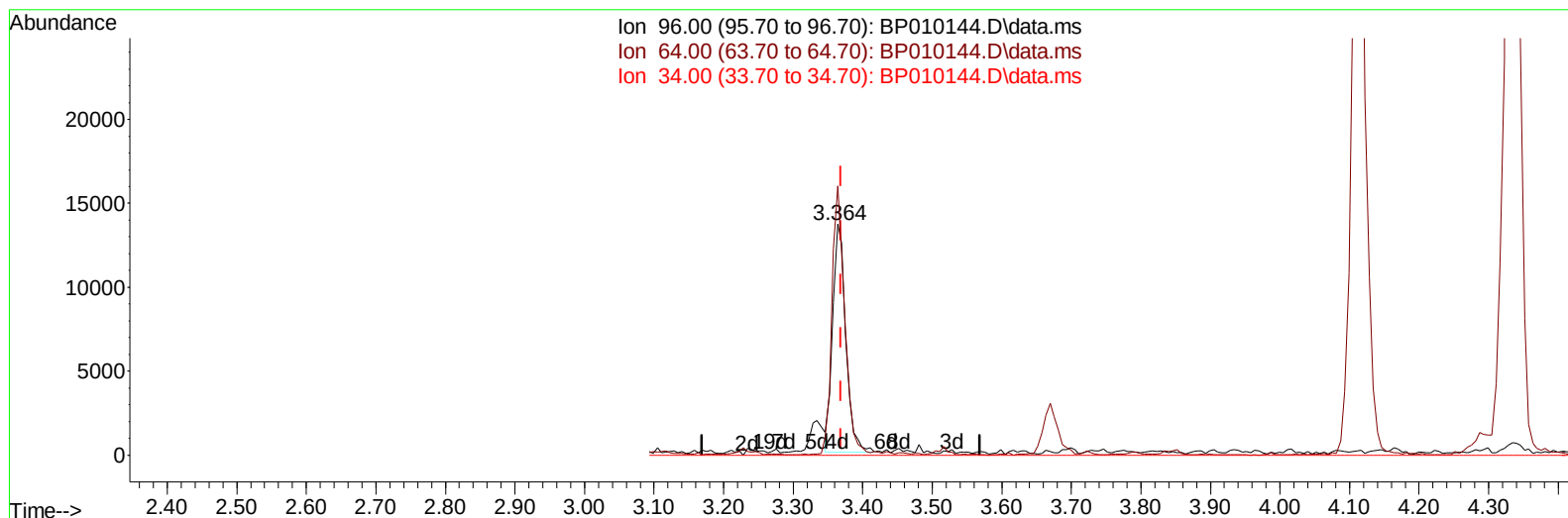
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TIC: BP010144.D\data.ms

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

3.364min (-0.005) 5.02 ng/uL m

response 18127

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	116.51#
34.00	0.00	0.00
0.00	0.00	0.00

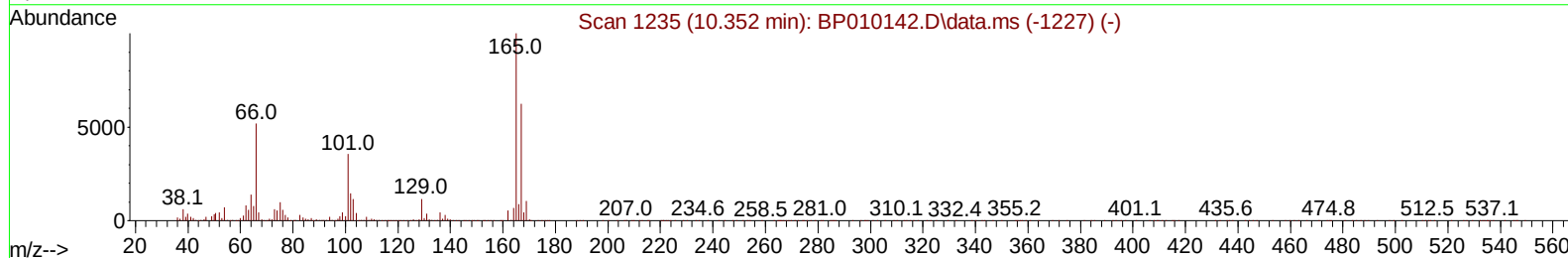
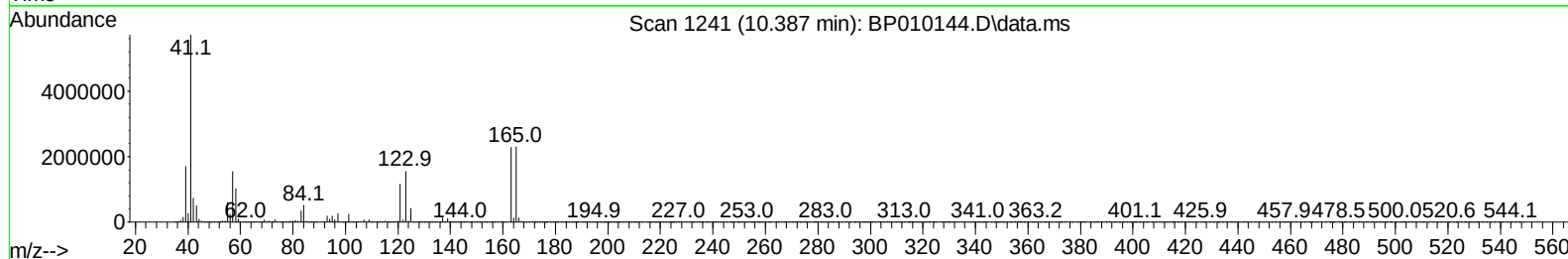
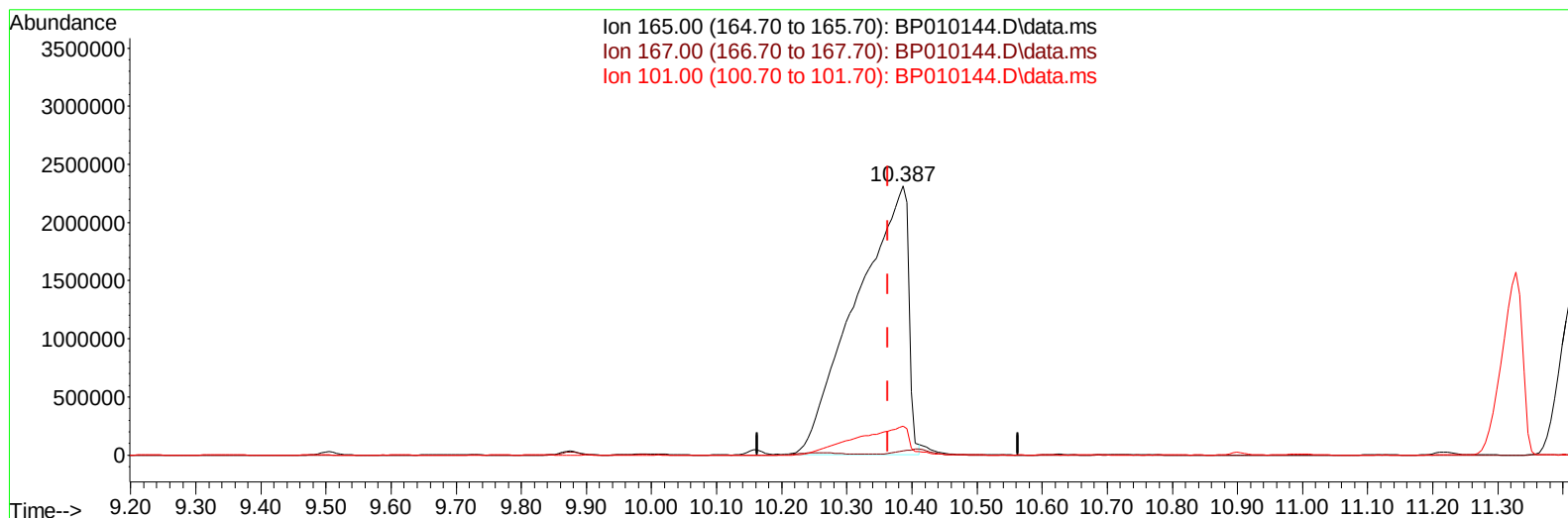
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Data File : BP010144.D
Acq On : 29 Apr 2022 11:03
Operator : CG/JU
Sample : N2587-02
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EXET0

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Response via : Initial Calibration



TIC: BP010144.D\data.ms

(28) 2,4-Dichlorophenol-d3 (S)

10.387min (+ 0.024) 1305.08 ng/uI

response 12711362

Ion	Exp%	Act%
165.00	100.00	100.00
167.00	57.70	1.49#
101.00	49.80	10.65#
0.00	0.00	0.00

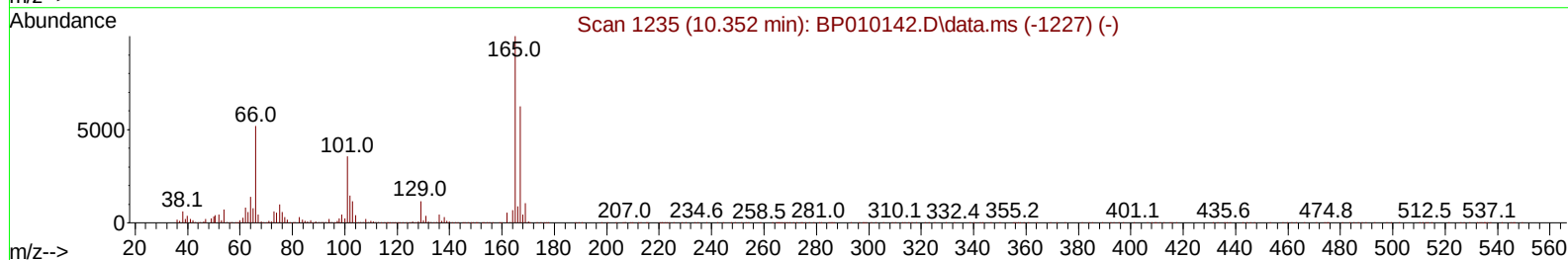
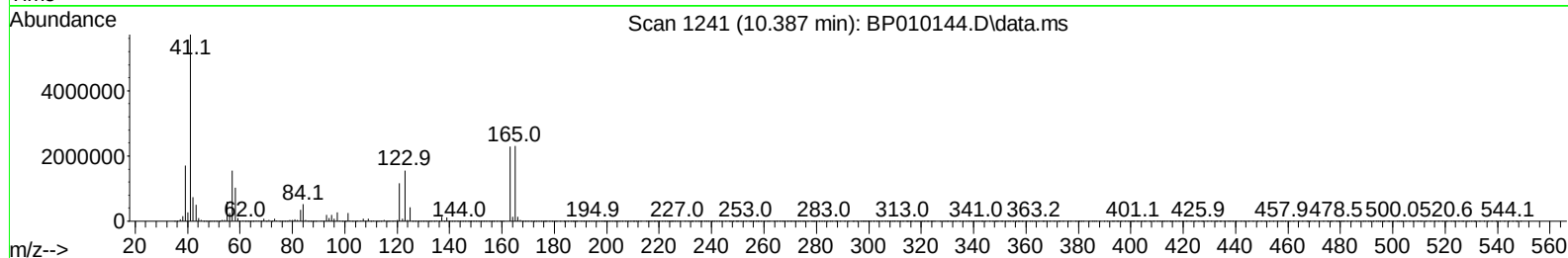
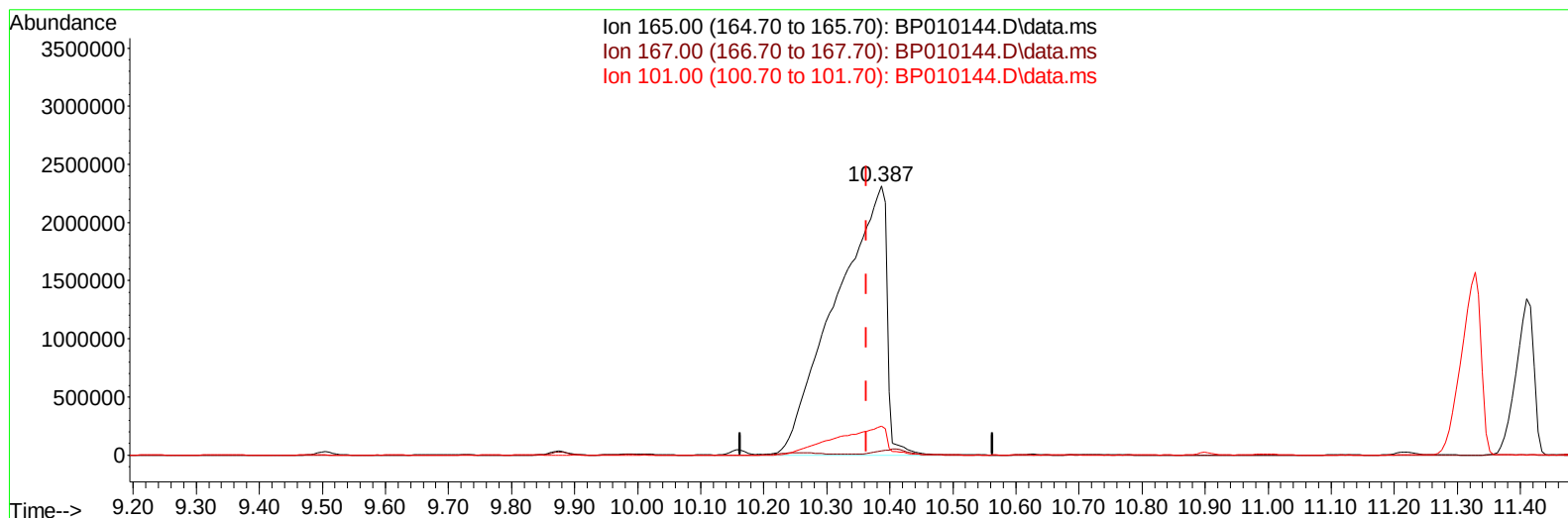
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Data File : BP010144.D
Acq On : 29 Apr 2022 11:03
Operator : CG/JU
Sample : N2587-02
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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TIC: BP010144.D\data.ms

(28) 2,4-Dichlorophenol-d3 (S)

10.387min (+ 0.024) 1324.40 ng/ul m

response 12899509

Ion	Exp%	Act%
165.00	100.00	100.00
167.00	57.70	1.49#
101.00	49.80	10.65#
0.00	0.00	0.00

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.934	152	157521	20.000	ng/ul	0.00
20) Naphthalene-d8	10.752	136	575986	20.000	ng/ul	# 0.00
38) Acenaphthene-d10	14.575	164	323523	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.328	188	630336	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.404	240	607454	20.000	ng/ul	0.00
88) Perylene-d12	23.857	264	650744	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.364	96	18127m	5.018	ng/uL	0.00
4) Pyridine-d5	3.787	84	73095	7.193	ng/ul	0.00
7) Phenol-d5	7.105	99	84513	5.898	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.258	67	259439	30.176	ng/ul	-0.01
11) 2-Chlorophenol-d4	7.463	132	255861	24.112	ng/ul	0.00
15) 4-Methylphenol-d8	8.652	113	172232	14.508	ng/ul	0.00
21) Nitrobenzene-d5	9.093	128	155405	33.923	ng/ul	0.00
24) 2-Nitrophenol-d4	9.834	143	161811	32.342	ng/ul	0.01
28) 2,4-Dichlorophenol-d3	10.387	165	12899509m	1324.398	ng/ul	0.02
31) 4-Chloroaniline-d4	10.910	131	205716	14.730	ng/ul	0.03
46) Dimethylphthalate-d6	13.993	166	810783	32.217	ng/ul	0.01
49) Acenaphthylene-d8	14.275	160	1015325	33.335	ng/ul	0.01
54) 4-Nitrophenol-d4	14.781	143	34719	7.494	ng/ul	0.02
60) Fluorene-d10	15.563	176	736399	34.016	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.675	200	140217	35.180	ng/ul	0.00
73) Anthracene-d10	17.422	188	1064324	34.902	ng/ul	0.00
81) Pyrene-d10	19.651	212	1211454	35.068	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.698	264	1181270	34.625	ng/ul	0.00
Target Compounds						
					Qvalue	
12) 2-Chlorophenol	7.499	128	33778	3.156	ng/ul	96
13) 2-Methylphenol	8.387	108	15870	1.437	ng/ul	100
30) Naphthalene	10.799	128	31329	1.039	ng/ul	99
78) Di-n-butylphthalate	18.275	149	50389	1.341	ng/ul#	97

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

