

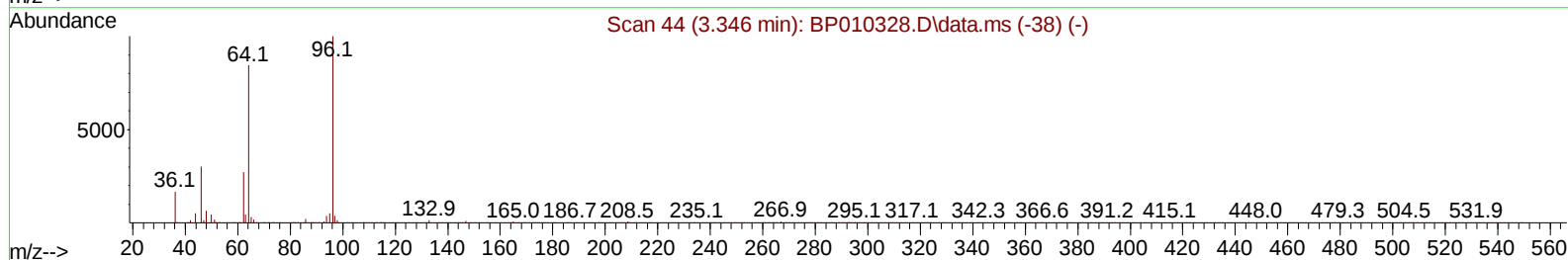
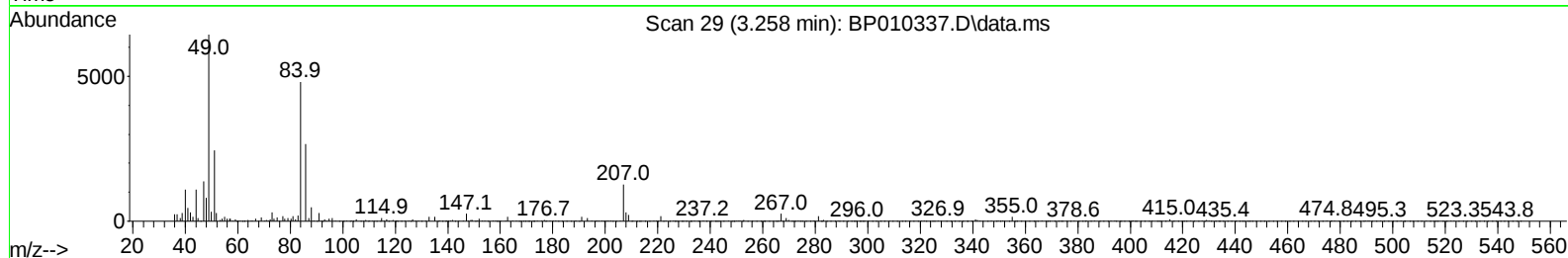
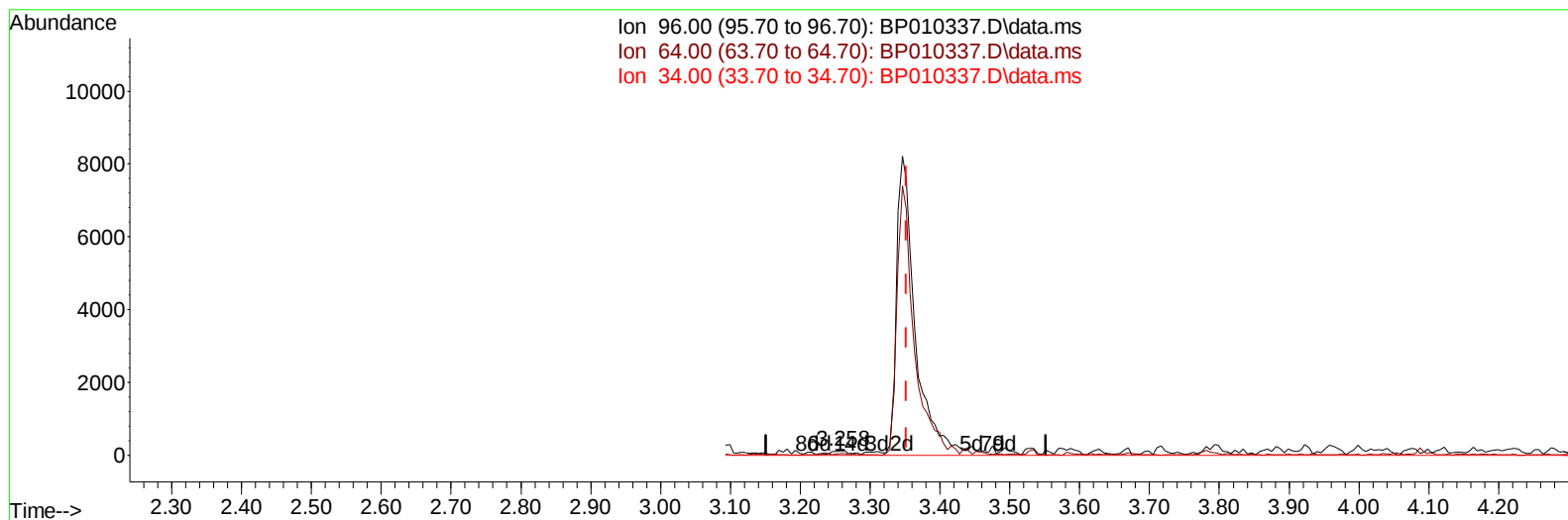
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051122\
 Data File : BP010337.D
 Acq On : 12 May 2022 23:05
 Operator : CG/JU
 Sample : N2801-01
 Misc :
 ALS Vial : 61 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 EXLB4

Manual IntegrationsAPPROVED

Quant Time: May 13 00:04:22 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP051022.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu May 12 01:10:15 2022
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 05/16/2022
 Supervised By :Yogesh Patel 05/17/2022



TIC: BP010337.D\data.ms

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

3.258min (-0.094) 0.05 ng/uL

response 157

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	44.92
34.00	0.00	0.00
0.00	0.00	0.00

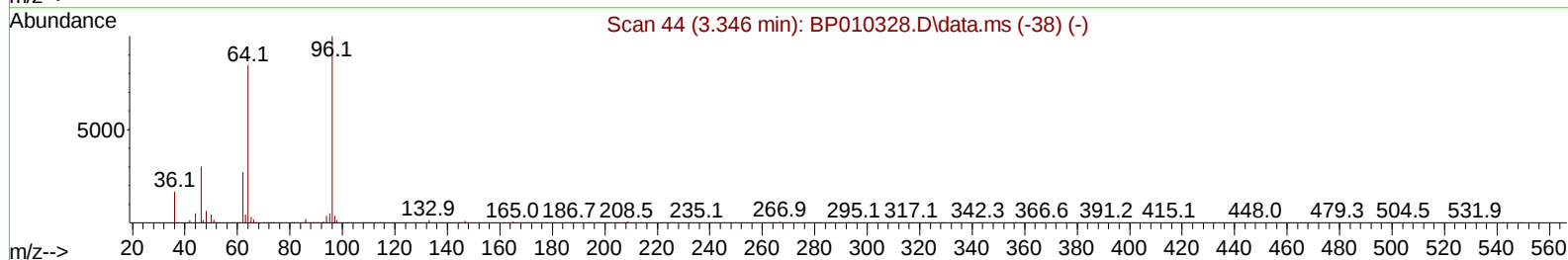
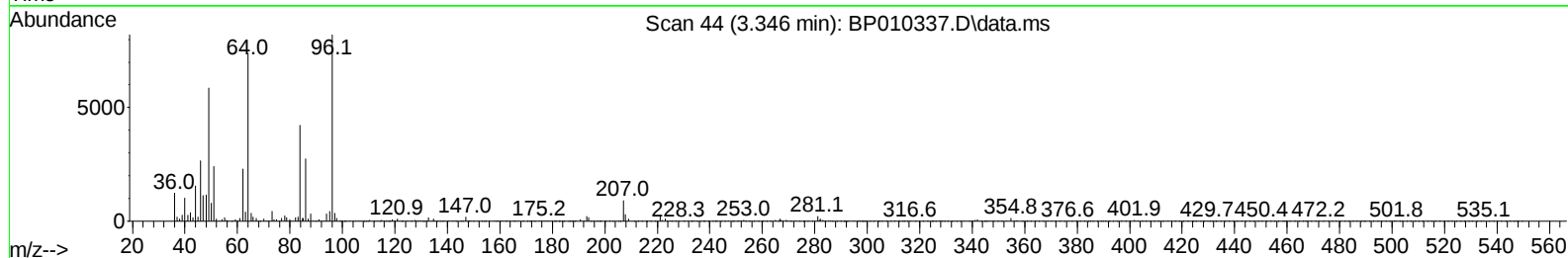
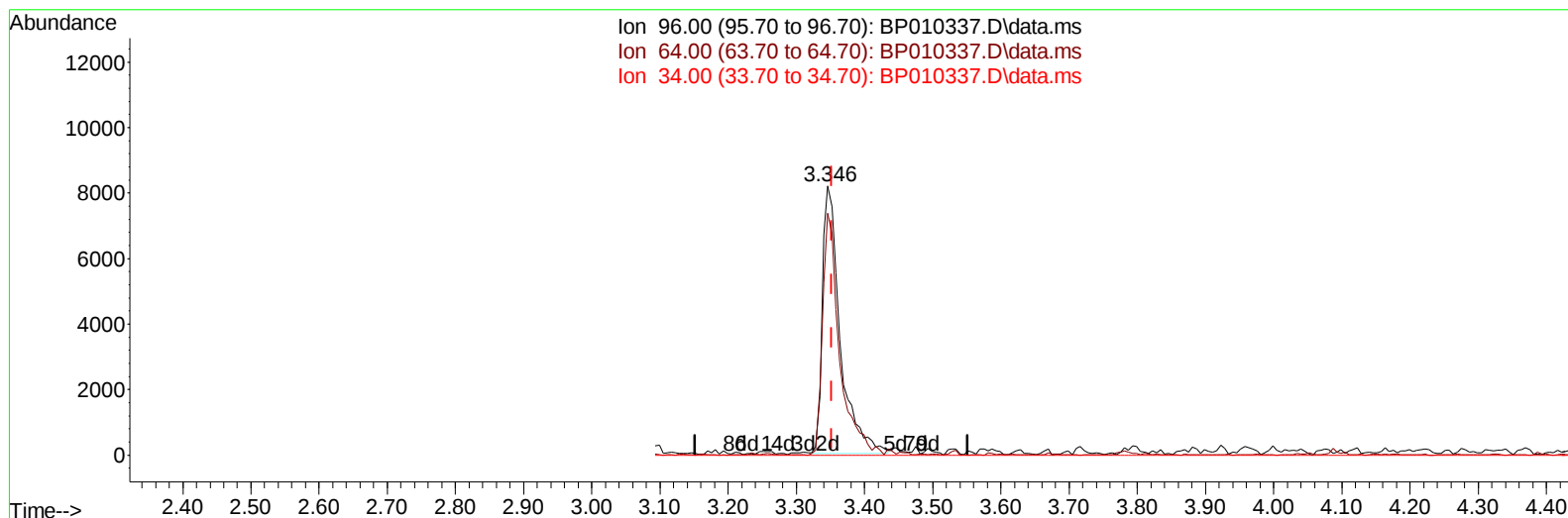
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TIC: BP010337.D\data.ms

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

3.346min (-0.006) 4.89 ng/uL m

response 14915

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	90.01#
34.00	0.00	0.00
0.00	0.00	0.00

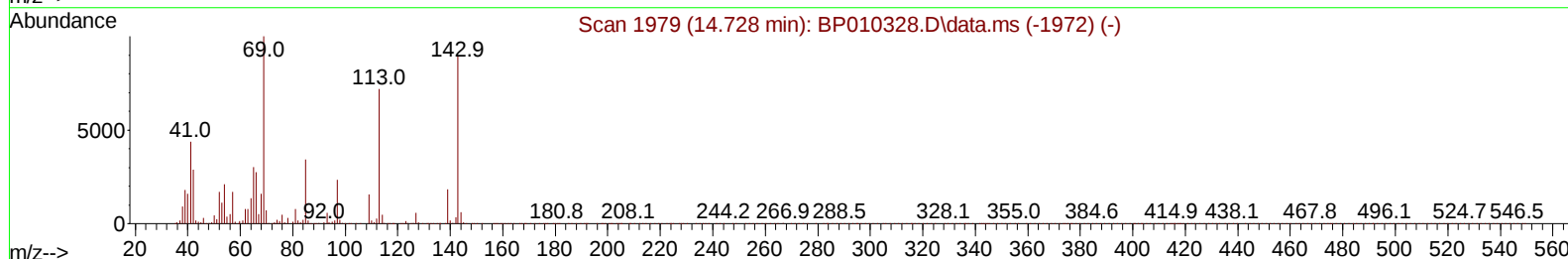
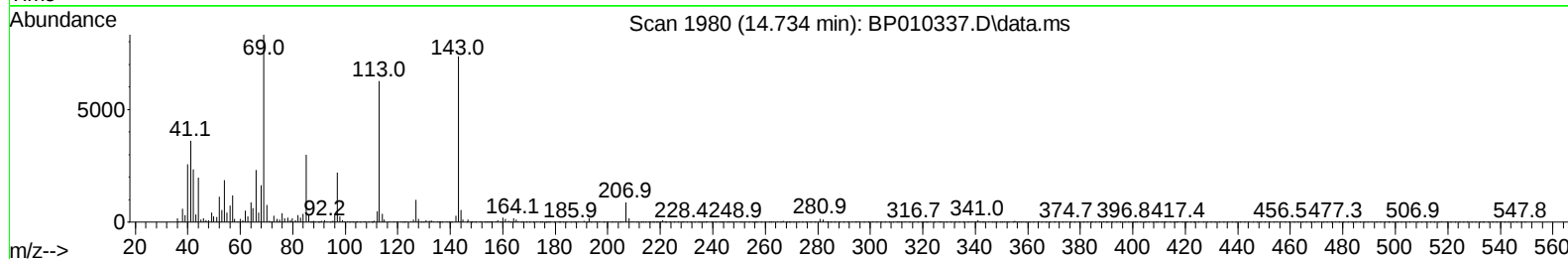
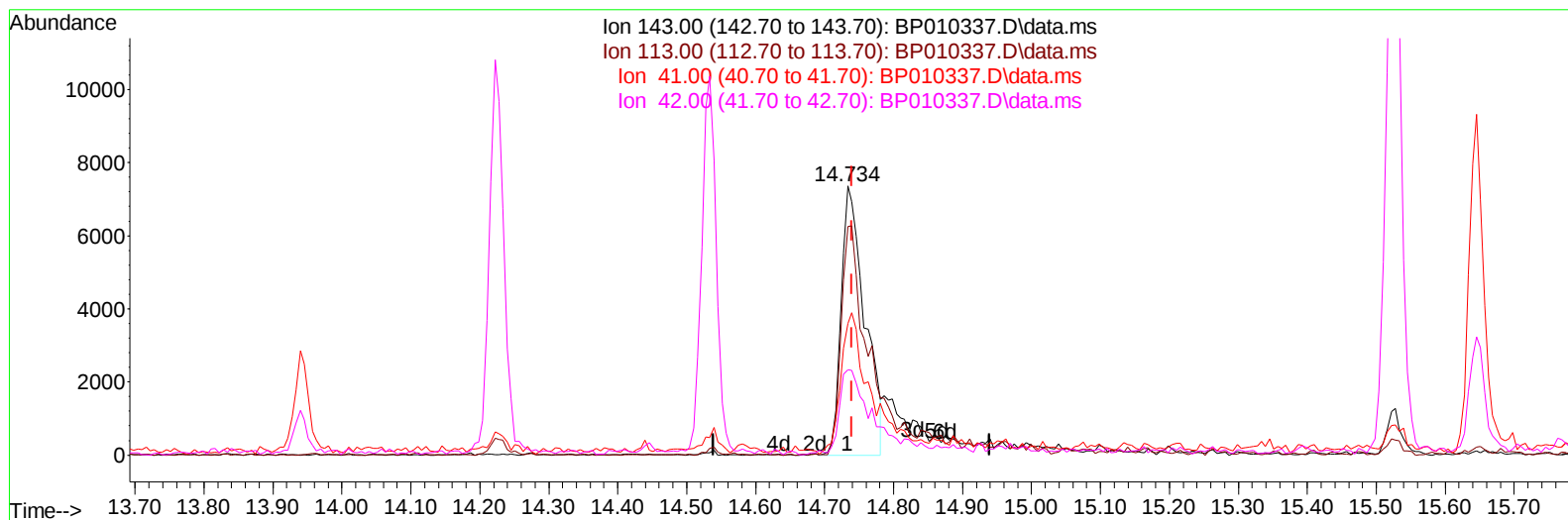
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Misc :
ALS Vial : 61 Sample Multiplier: 1

Instrument :
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ClientSampleId :
EXLB4

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TIC: BP010337.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.734min (-0.006) 2.77 ng/u1

response 17437

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	132.70	85.09#
41.00	23.20	48.85#
42.00	18.00	31.81#

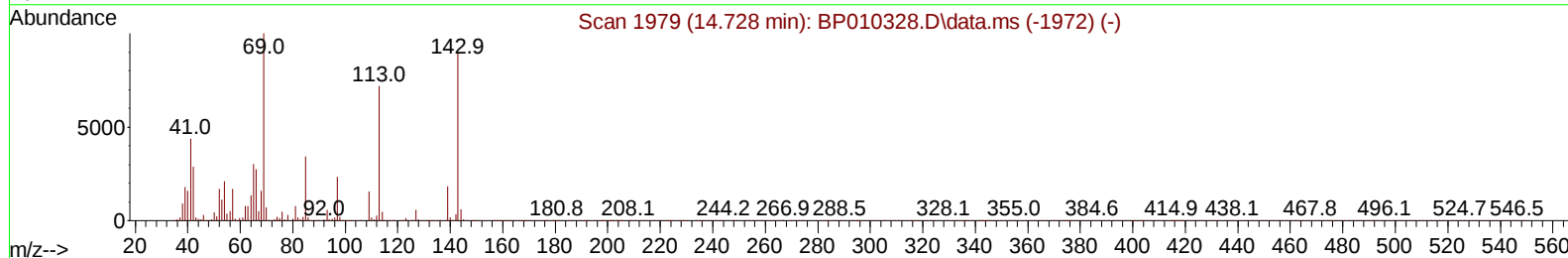
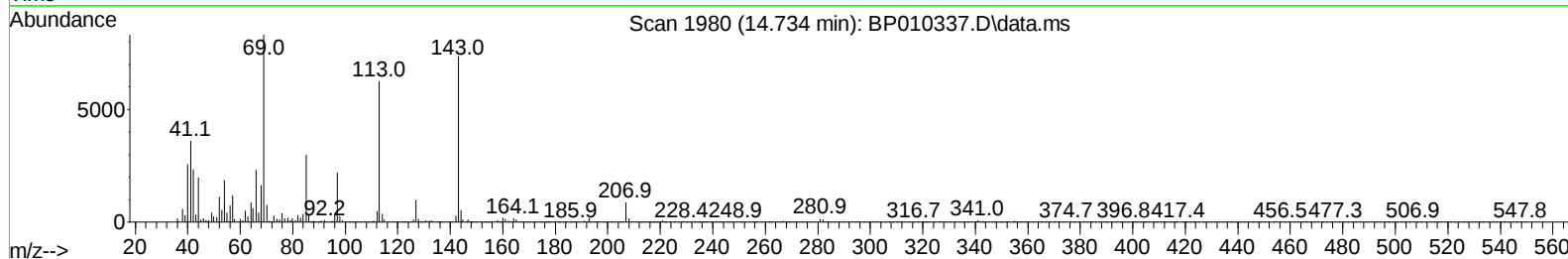
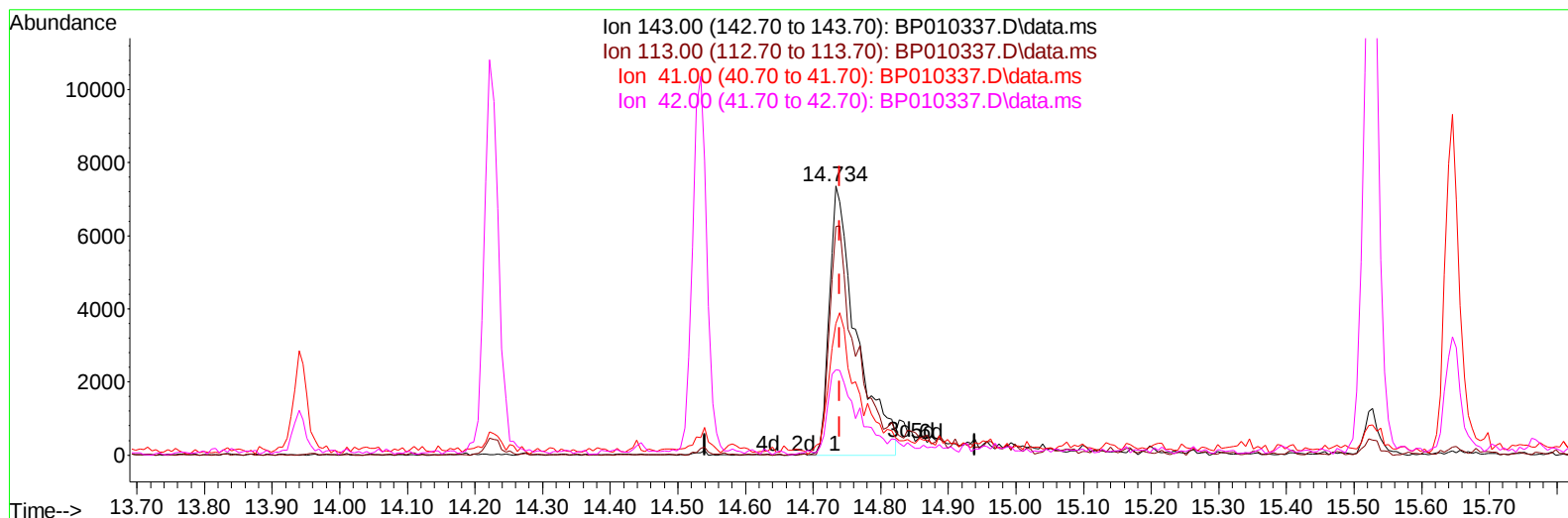
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Operator : CG/JU
Sample : N2801-01
Misc :
ALS Vial : 61 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EXLB4

Manual IntegrationsAPPROVED

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TIC: BP010337.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.734min (-0.006) 3.24 ng/ul m

response 20394

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	132.70	85.09#
41.00	23.20	48.85#
42.00	18.00	31.81#

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 Sample : N2801-01
 Misc :
 ALS Vial : 61 Sample Multiplier: 1

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.904	152	129485	20.000	ng/ul	0.00
20) Naphthalene-d8	10.704	136	574994	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.534	164	412435	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.286	188	905024	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.363	240	939157	20.000	ng/ul	#-0.01
88) Perylene-d12	23.792	264	910835	20.000	ng/ul	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.346	96	14915m	4.893	ng/uL	0.00
4) Pyridine-d5	3.775	84	52872	6.234	ng/ul	0.00
7) Phenol-d5	7.069	99	61760	5.135	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.228	67	220363	29.619	ng/ul	0.00
11) 2-Chlorophenol-d4	7.434	132	187865	21.211	ng/ul	0.00
15) 4-Methylphenol-d8	8.610	113	126685	12.813	ng/ul	0.00
21) Nitrobenzene-d5	9.063	128	132771	28.798	ng/ul	0.00
24) 2-Nitrophenol-d4	9.787	143	140755	26.777	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.328	165	234739	24.535	ng/ul	0.00
31) 4-Chloroaniline-d4	10.840	131	307139	21.745	ng/ul	0.00
46) Dimethylphthalate-d6	13.945	166	968805	29.485	ng/ul	0.00
49) Acenaphthylene-d8	14.228	160	1083888	27.668	ng/ul	0.00
54) 4-Nitrophenol-d4	14.734	143	20394m	3.236	ng/ul	0.00
60) Fluorene-d10	15.528	176	850499	30.665	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.645	200	139134	22.450	ng/ul	-0.01
73) Anthracene-d10	17.386	188	1428169	32.225	ng/ul	0.00
81) Pyrene-d10	19.616	212	1767002	34.047	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.639	264	1632014	33.598	ng/ul	-0.02

Target Compounds Qvalue

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

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