

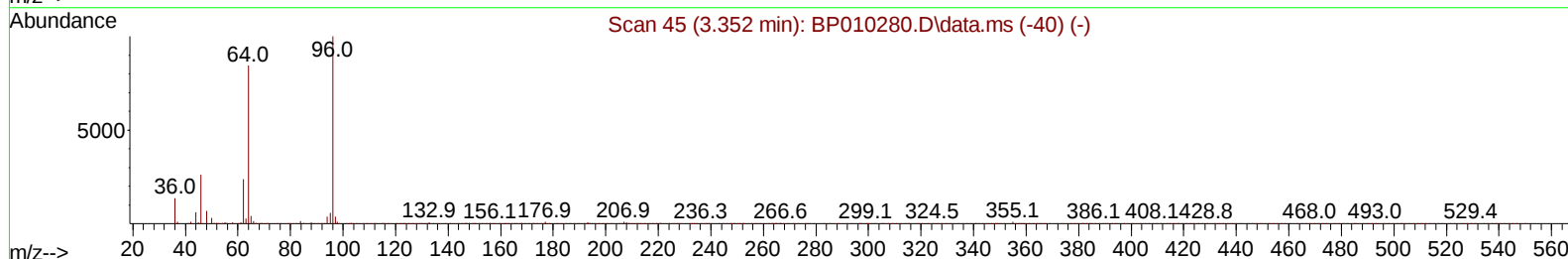
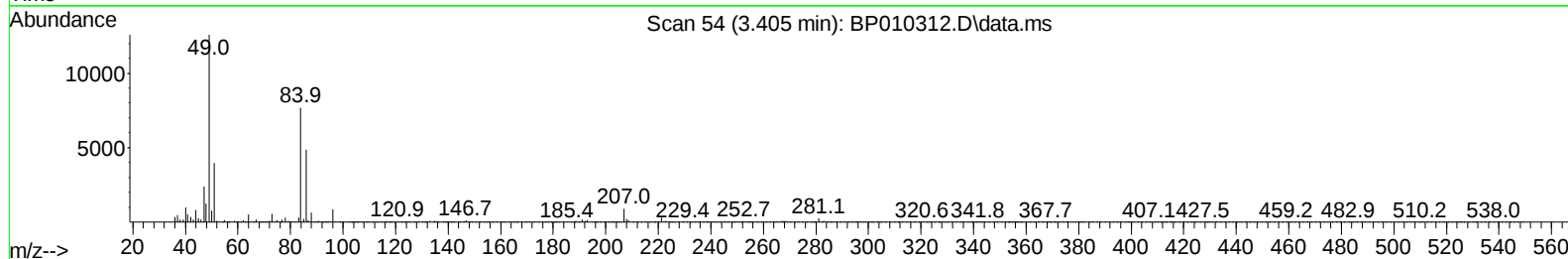
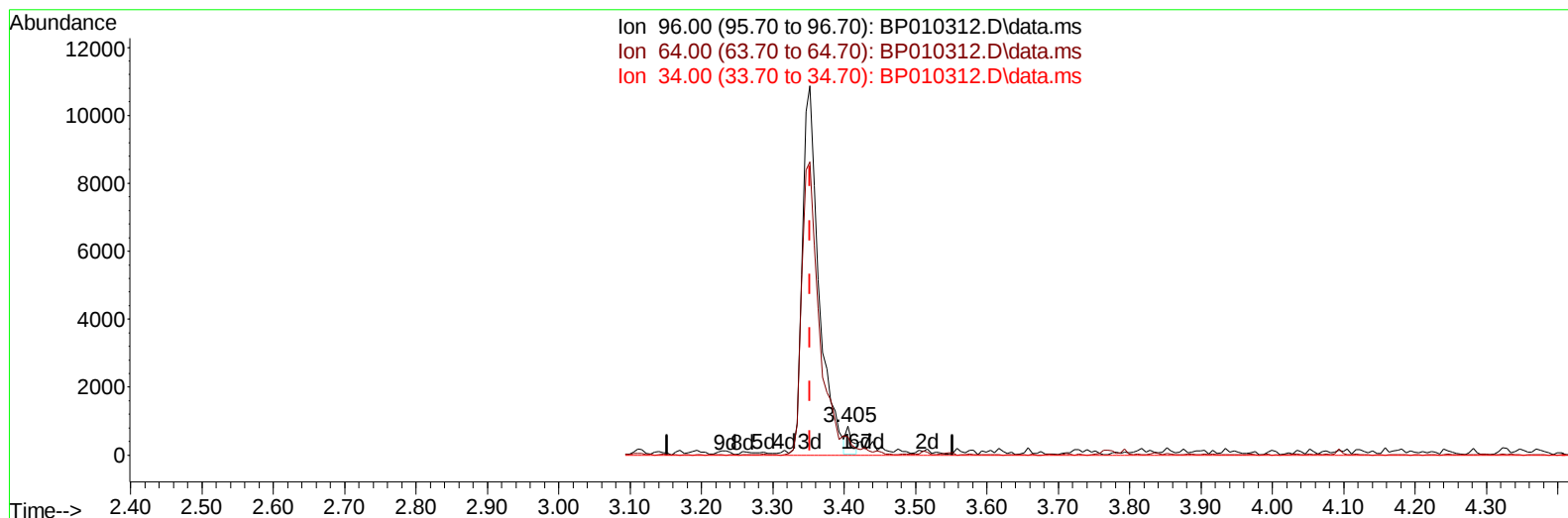
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051122\
 Data File : BP010312.D
 Acq On : 12 May 2022 07:12
 Operator : CG/JU
 Sample : N2668-18
 Misc :
 ALS Vial : 36 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 EXL76

Manual IntegrationsAPPROVED

Quant Time: May 12 08:22:27 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 05/12/2022
 Supervised By :Yogesh Patel 05/17/2022



TIC: BP010312.D\data.ms

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

3.405min (+ 0.053) 0.13 ng/uL

response 516

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	63.31
34.00	0.00	0.00
0.00	0.00	0.00

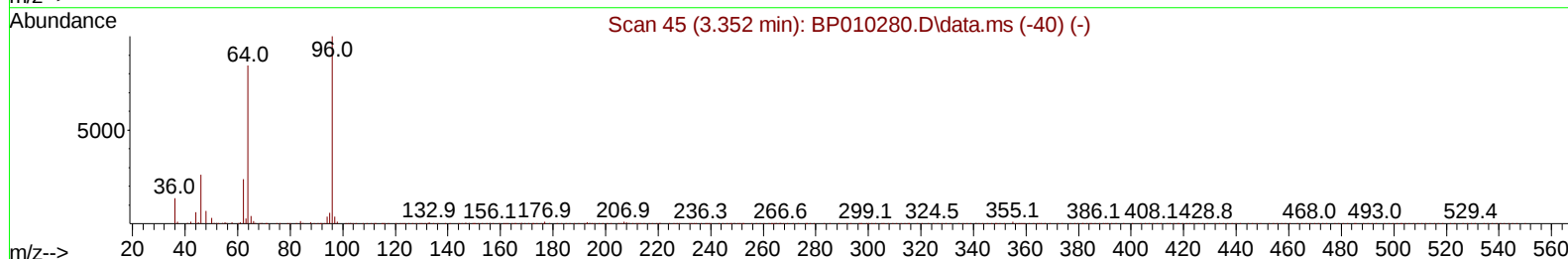
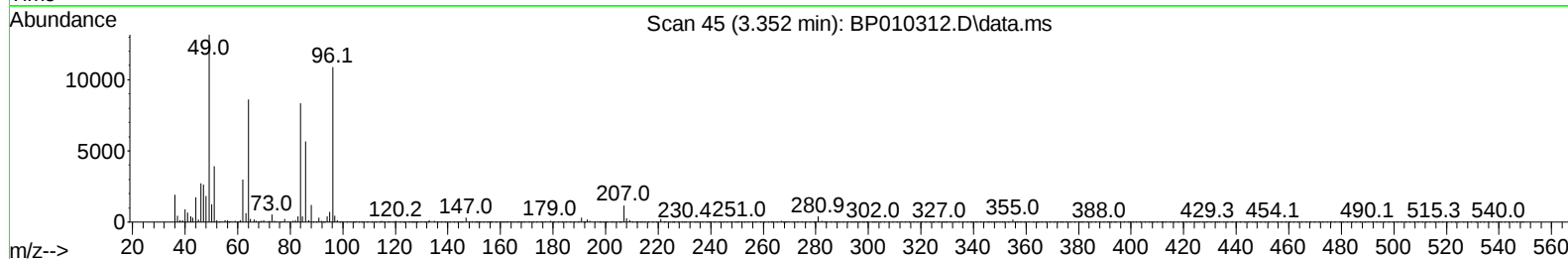
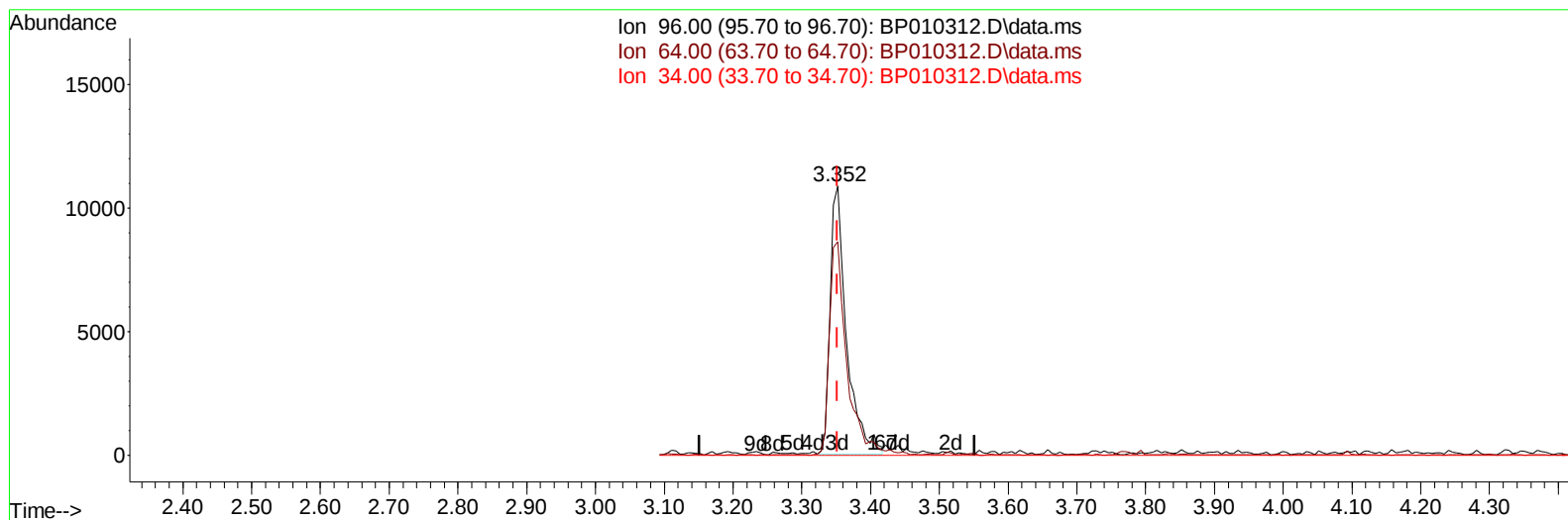
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TIC: BP010312.D\data.ms

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

3.352min (+ 0.000) 4.57 ng/uL m

response 17965

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	79.36#
34.00	0.00	0.00
0.00	0.00	0.00

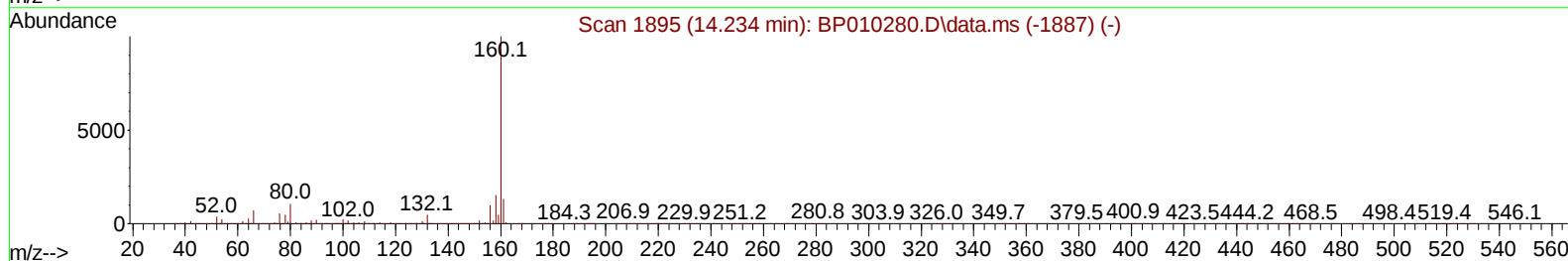
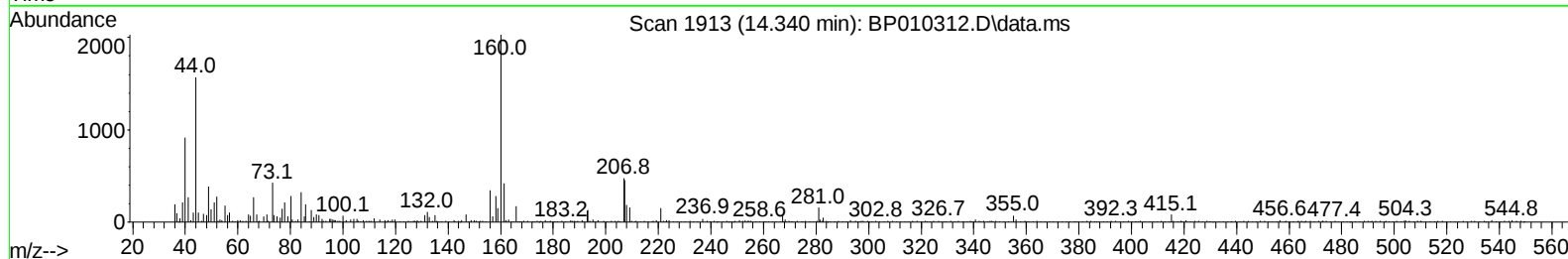
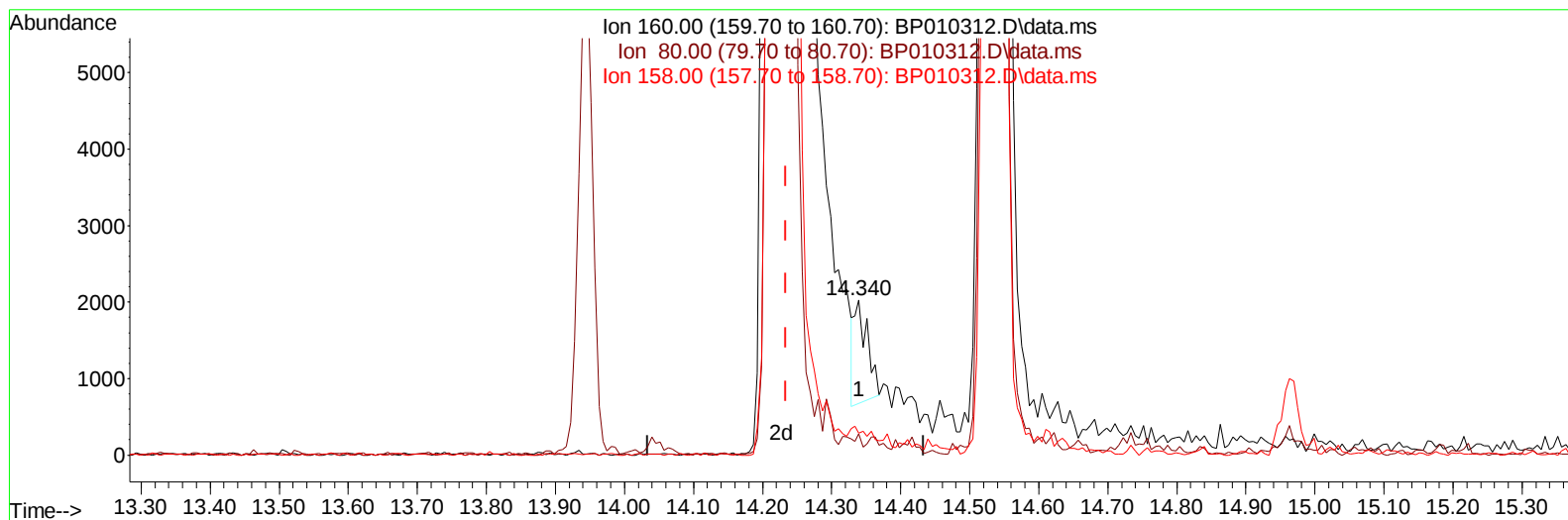
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Misc :
ALS Vial : 36 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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TIC: BP010312.D\data.ms

(49) Acenaphthylene-d8 (S)

14.340min (+ 0.106) 0.04 ng/u1

response 1800

Ion	Exp%	Act%
160.00	100.00	100.00
80.00	17.00	13.74
158.00	13.10	13.99
0.00	0.00	0.00

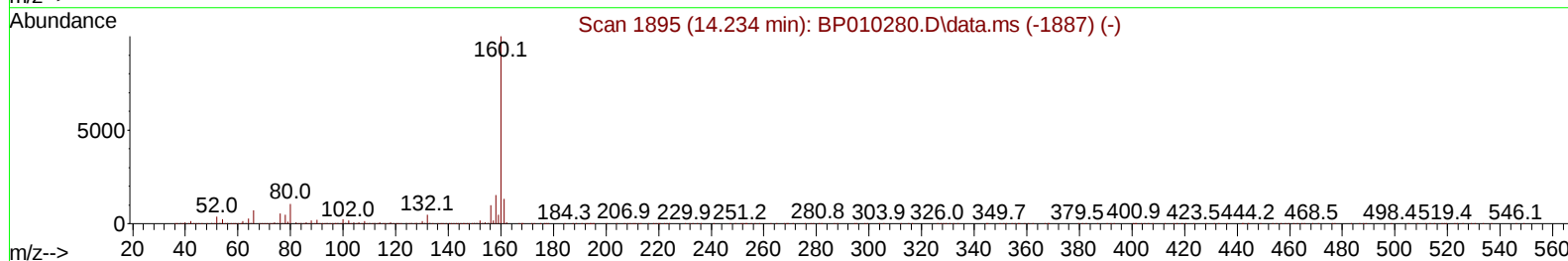
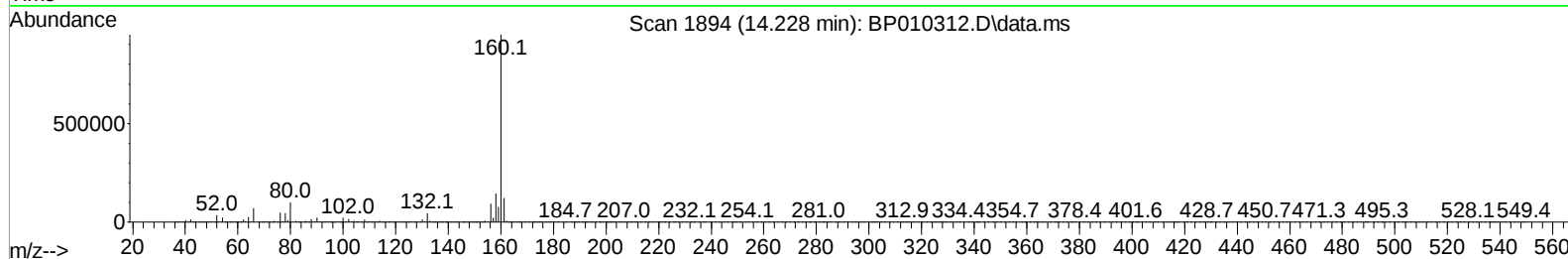
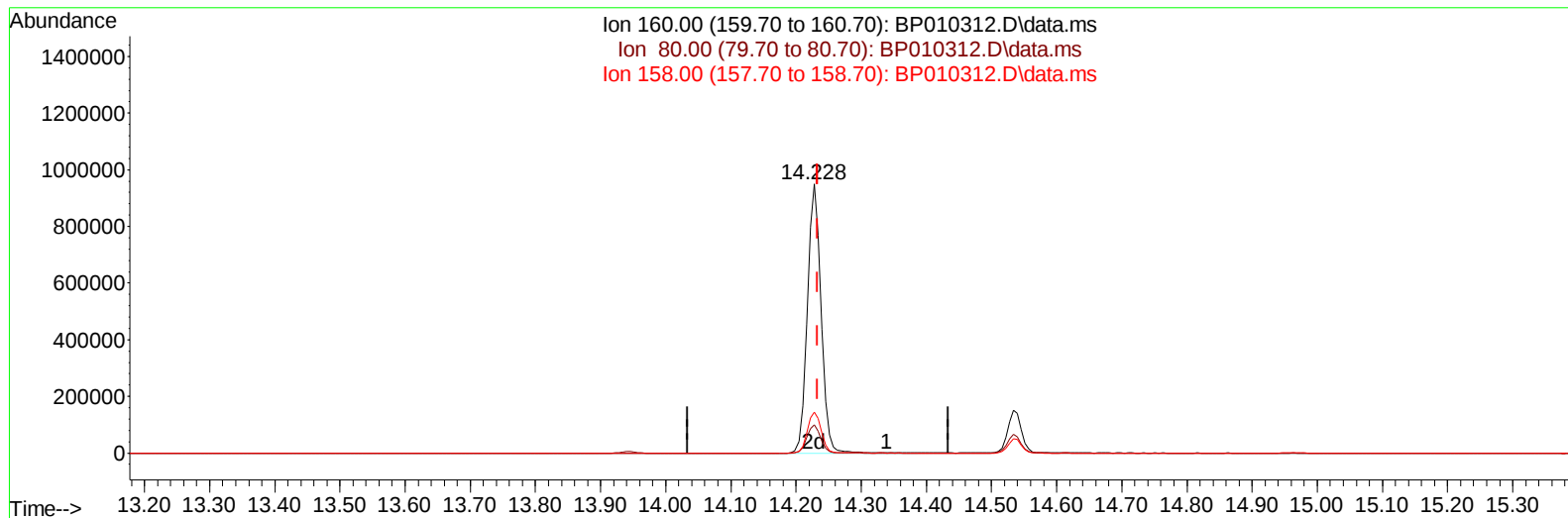
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Operator : CG/JU
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Misc :
ALS Vial : 36 Sample Multiplier: 1

Instrument :
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TIC: BP010312.D\data.ms

(49) Acenaphthylene-d8 (S)

14.228min (-0.006) 29.11 ng/ul m

response 1397590

Ion	Exp%	Act%
160.00	100.00	100.00
80.00	17.00	10.45#
158.00	13.10	15.31
0.00	0.00	0.00

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 Misc :
 ALS Vial : 36 Sample Multiplier: 1

Instrument :
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Manual IntegrationsAPPROVED

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.905	152	166975	20.000	ng/ul	0.00
20) Naphthalene-d8	10.704	136	723121	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.534	164	505392	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.286	188	1109160	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.369	240	1107047	20.000	ng/ul	# 0.00
88) Perylene-d12	23.798	264	1059998	20.000	ng/ul	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.352	96	17965m	4.571	ng/uL	0.00
4) Pyridine-d5	3.775	84	82716	7.562	ng/ul	0.00
7) Phenol-d5	7.069	99	76252	4.917	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.234	67	287657	29.983	ng/ul	0.00
11) 2-Chlorophenol-d4	7.434	132	247185	21.642	ng/ul	0.00
15) 4-Methylphenol-d8	8.610	113	160553	12.592	ng/ul	0.00
21) Nitrobenzene-d5	9.063	128	176809	30.494	ng/ul	0.00
24) 2-Nitrophenol-d4	9.787	143	185635	28.081	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.328	165	313692	26.071	ng/ul	0.00
31) 4-Chloroaniline-d4	10.840	131	414271	23.322	ng/ul	0.00
46) Dimethylphthalate-d6	13.945	166	1289877	32.036	ng/ul	0.00
49) Acenaphthylene-d8	14.228	160	1397590m	29.114	ng/ul	0.00
54) 4-Nitrophenol-d4	14.734	143	27784	3.598	ng/ul	0.00
60) Fluorene-d10	15.528	176	1097747	32.299	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.645	200	209494	27.582	ng/ul	-0.01
73) Anthracene-d10	17.386	188	1903675	35.048	ng/ul	0.00
81) Pyrene-d10	19.616	212	2314081	37.826	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.645	264	2066166	36.550	ng/ul	-0.01

Target Compounds Qvalue

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

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