

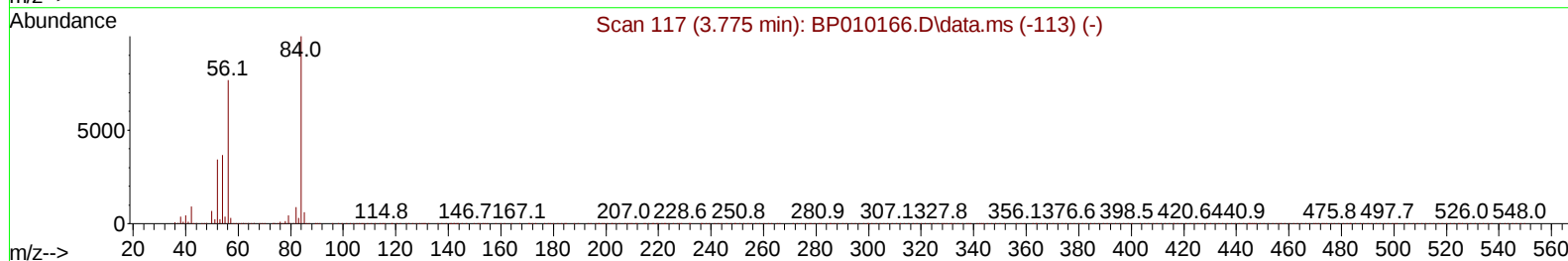
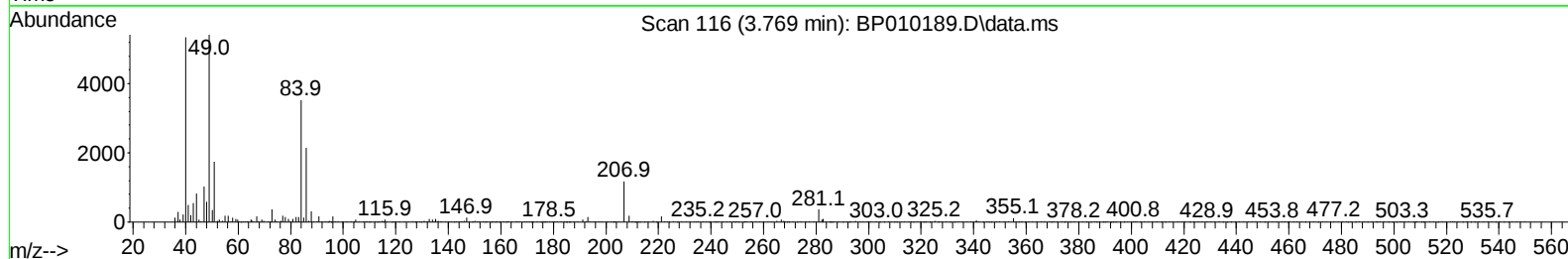
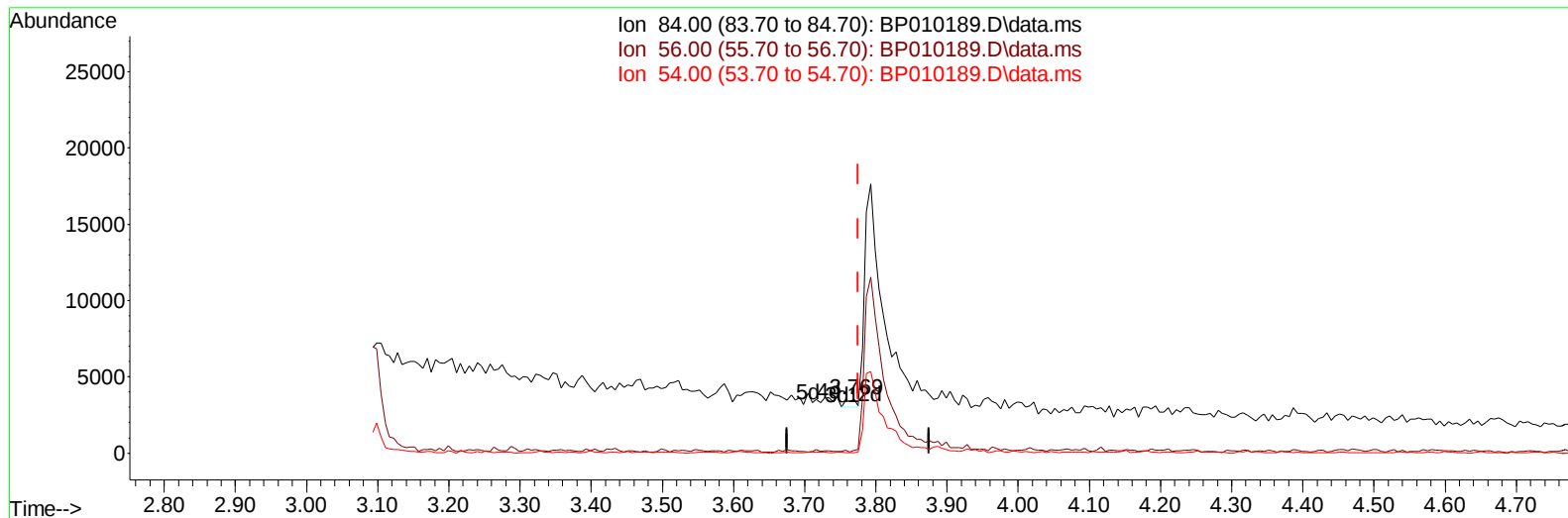
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP050222\
 Data File : BP010189.D
 Acq On : 03 May 2022 00:50
 Operator : CG/JU
 Sample : N2615-08
 Misc :
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BH0A3

Manual IntegrationsAPPROVED

Quant Time: May 03 02:53:40 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP042122.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue May 03 00:58:32 2022
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 05/03/2022
 Supervised By :Yogesh Patel 05/05/2022



TIC: BP010189.D\data.ms

(4) Pyridine-d5 (S)

3.769min (-0.006) 0.05 ng/u1

response 492

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	30.50	5.06#
54.00	24.80	0.85#
0.00	0.00	0.00

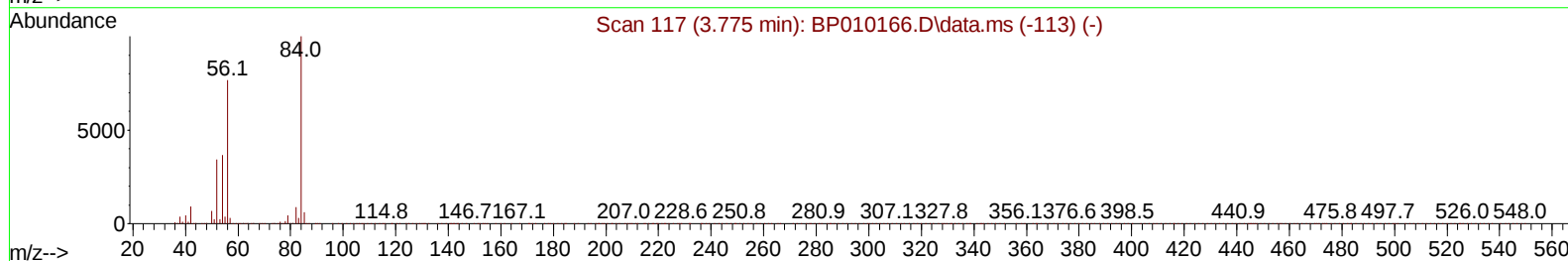
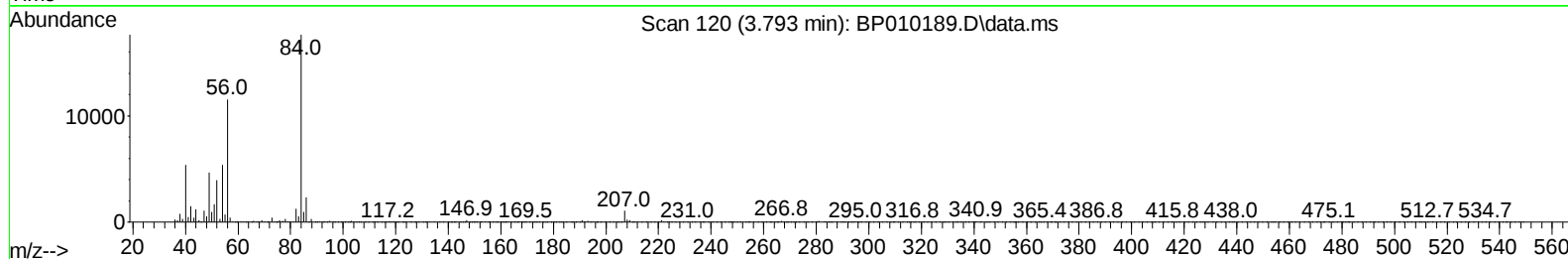
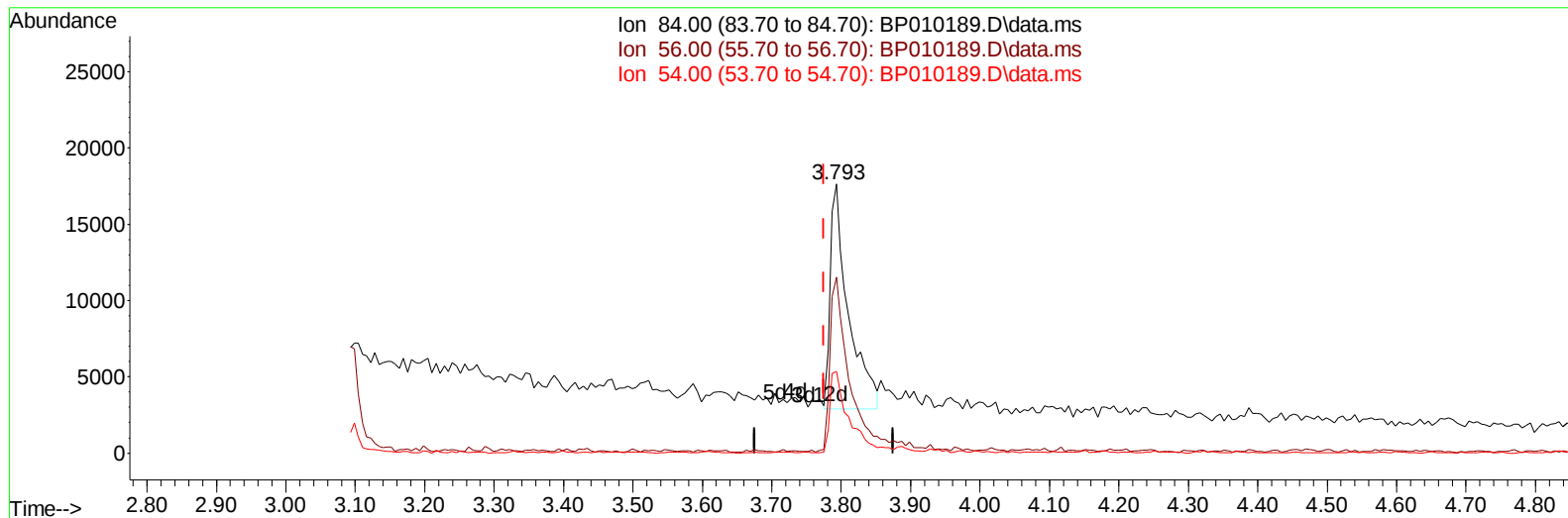
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TIC: BP010189.D\data.ms

(4) Pyridine-d5 (S)

3.793min (+ 0.017) 2.74 ng/u1 m

response 26789

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	30.50	65.43#
54.00	24.80	30.42#
0.00	0.00	0.00

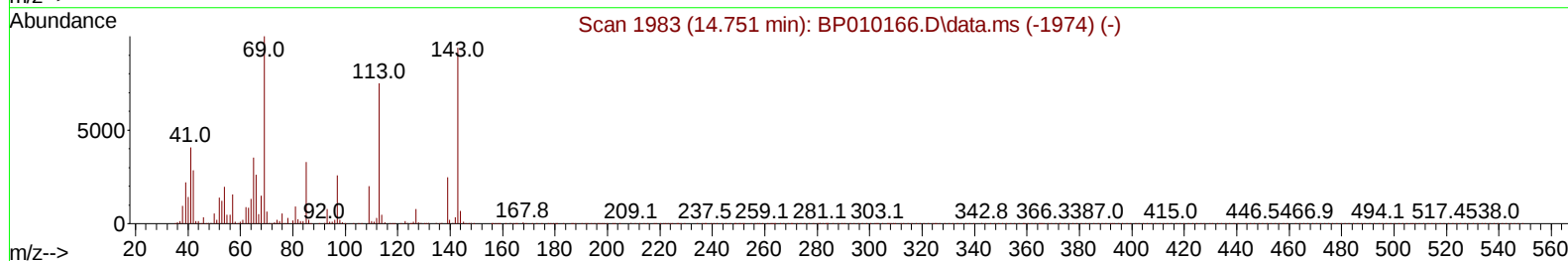
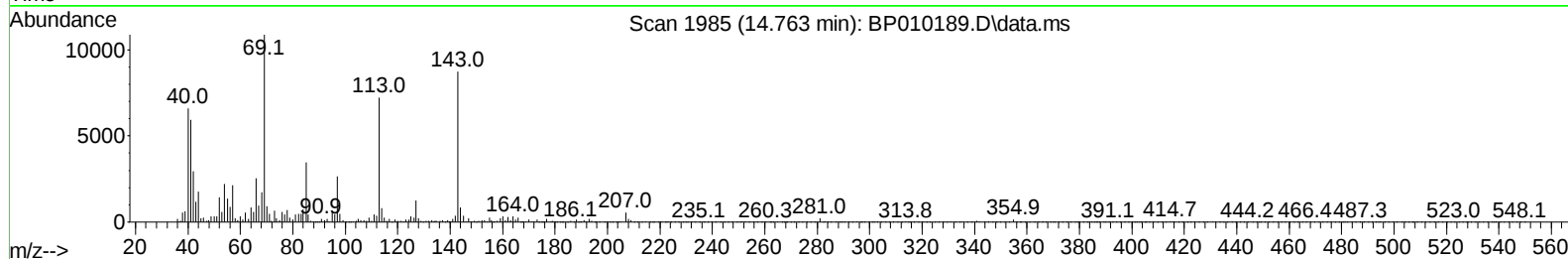
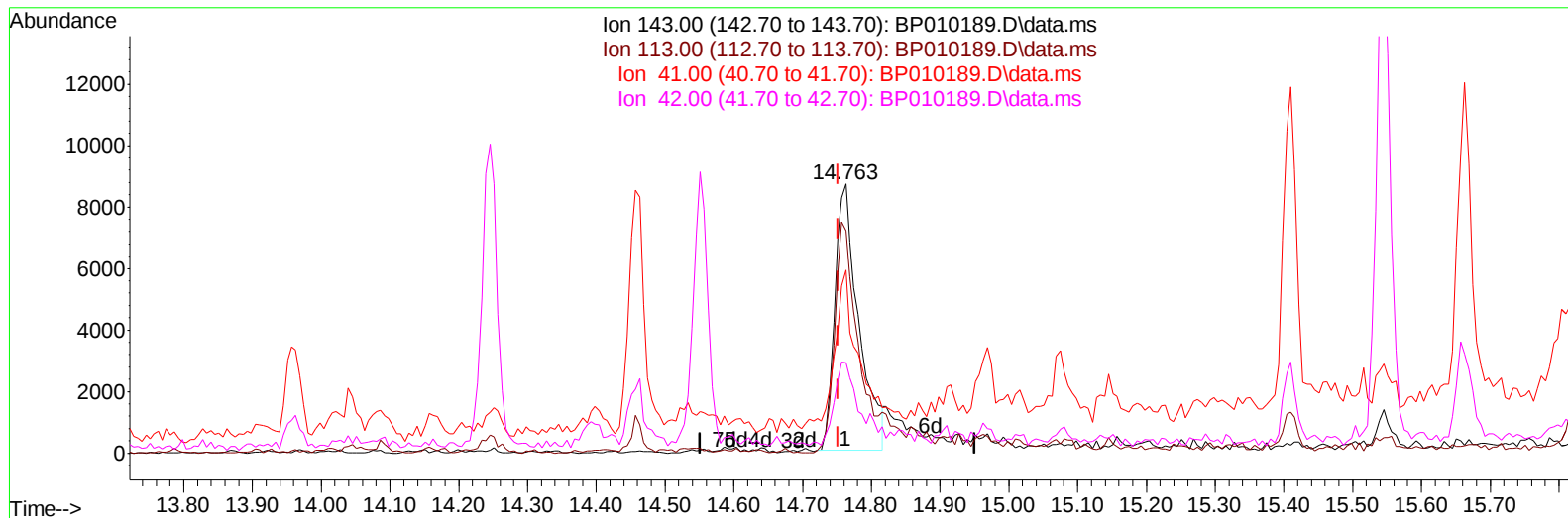
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Instrument :
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TIC: BP010189.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.763min (+ 0.012) 3.89 ng/u1

response 20906

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	132.70	82.64#
41.00	23.20	68.01#
42.00	18.00	33.59#

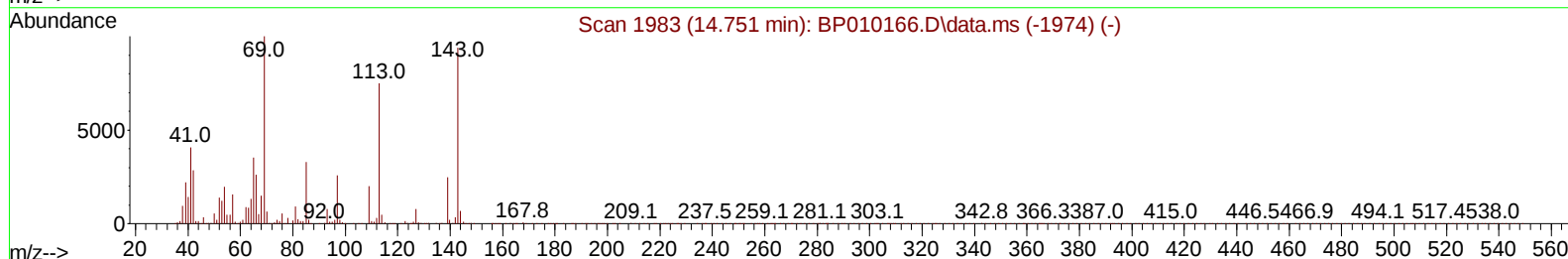
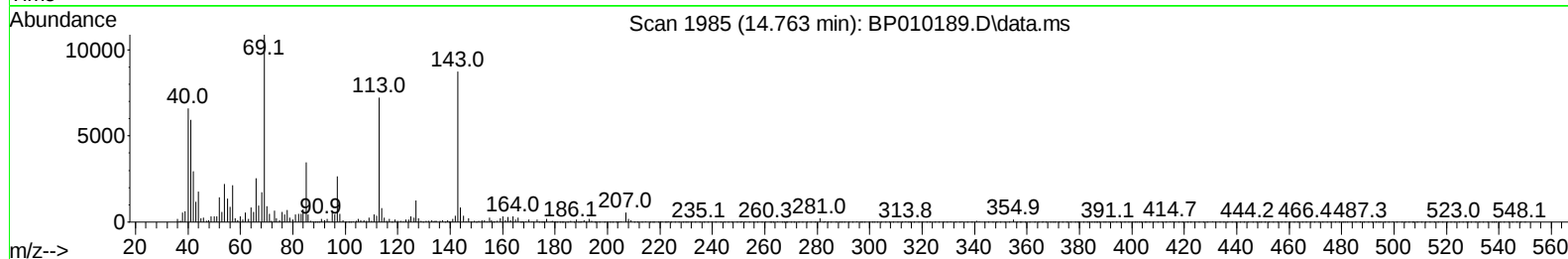
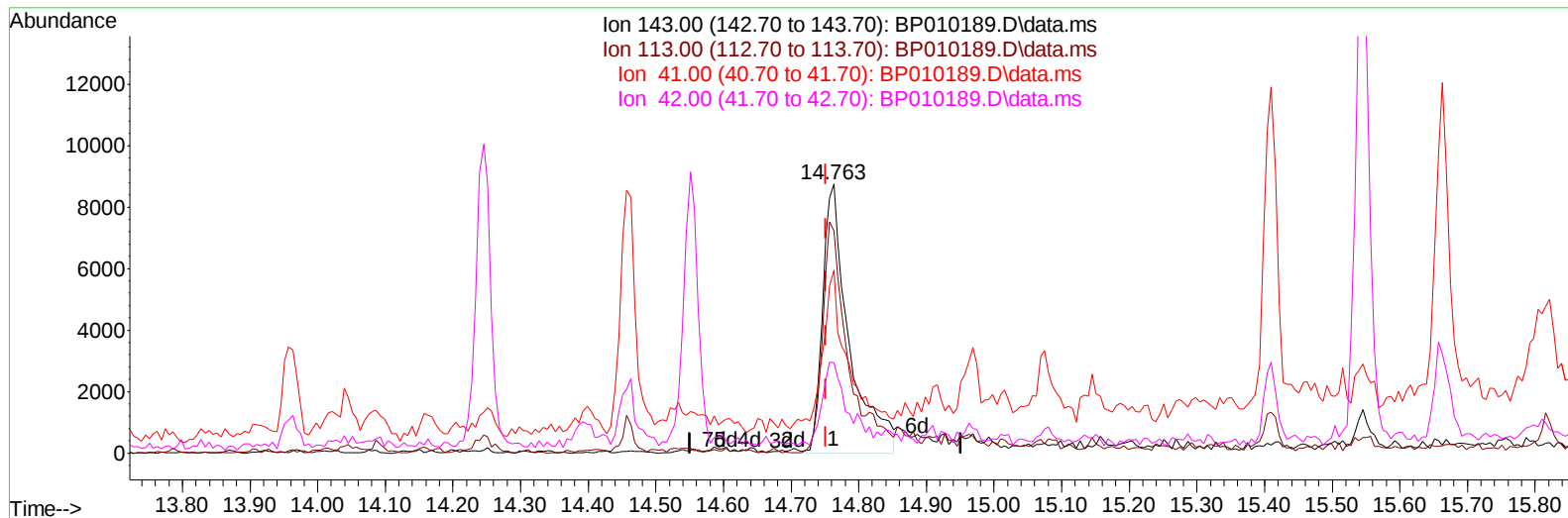
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 Operator : CG/JU
 Sample : N2615-08
 Misc :
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
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Manual IntegrationsAPPROVED

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TIC: BP010189.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.763min (+ 0.012) 4.42 ng/ul m

response 23773

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	132.70	82.64#
41.00	23.20	68.01#
42.00	18.00	33.59#

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 Misc :
 ALS Vial : 25 Sample Multiplier: 1

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.922	152	151326	20.000	ng/ul	0.00
20) Naphthalene-d8	10.728	136	622136	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.551	164	375313	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.304	188	720347	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.386	240	674147	20.000	ng/ul	# 0.00
88) Perylene-d12	23.827	264	669612	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.358	96	15426	4.445	ng/uL	0.00
4) Pyridine-d5	3.793	84	26789m	2.744	ng/ul	0.02
7) Phenol-d5	7.087	99	77401	5.623	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.246	67	278207	33.684	ng/ul	0.00
11) 2-Chlorophenol-d4	7.451	132	254006	24.918	ng/ul	0.00
15) 4-Methylphenol-d8	8.628	113	158633	13.910	ng/ul	0.00
21) Nitrobenzene-d5	9.081	128	166817	33.713	ng/ul	0.00
24) 2-Nitrophenol-d4	9.804	143	178508	33.032	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.345	165	277361	26.364	ng/ul	0.00
31) 4-Chloroaniline-d4	10.863	131	267334	17.722	ng/ul	0.00
46) Dimethylphthalate-d6	13.963	166	1055950	36.169	ng/ul	0.00
49) Acenaphthylene-d8	14.245	160	1156662	32.735	ng/ul	0.00
54) 4-Nitrophenol-d4	14.763	143	23773m	4.423	ng/ul	0.01
60) Fluorene-d10	15.545	176	883792	35.191	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.663	200	152302	33.437	ng/ul	0.00
73) Anthracene-d10	17.404	188	1419129	40.721	ng/ul	0.00
81) Pyrene-d10	19.633	212	1632896	42.591	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.674	264	1481670	42.207	ng/ul	0.00

Target Compounds Qvalue

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

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