

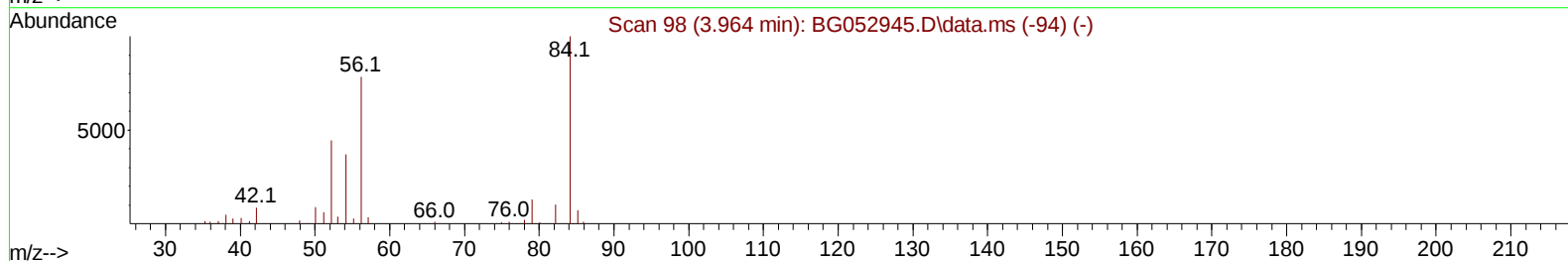
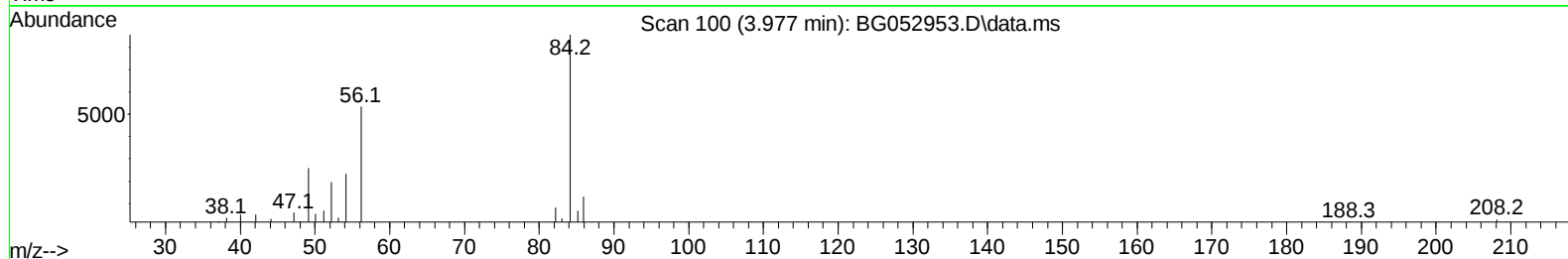
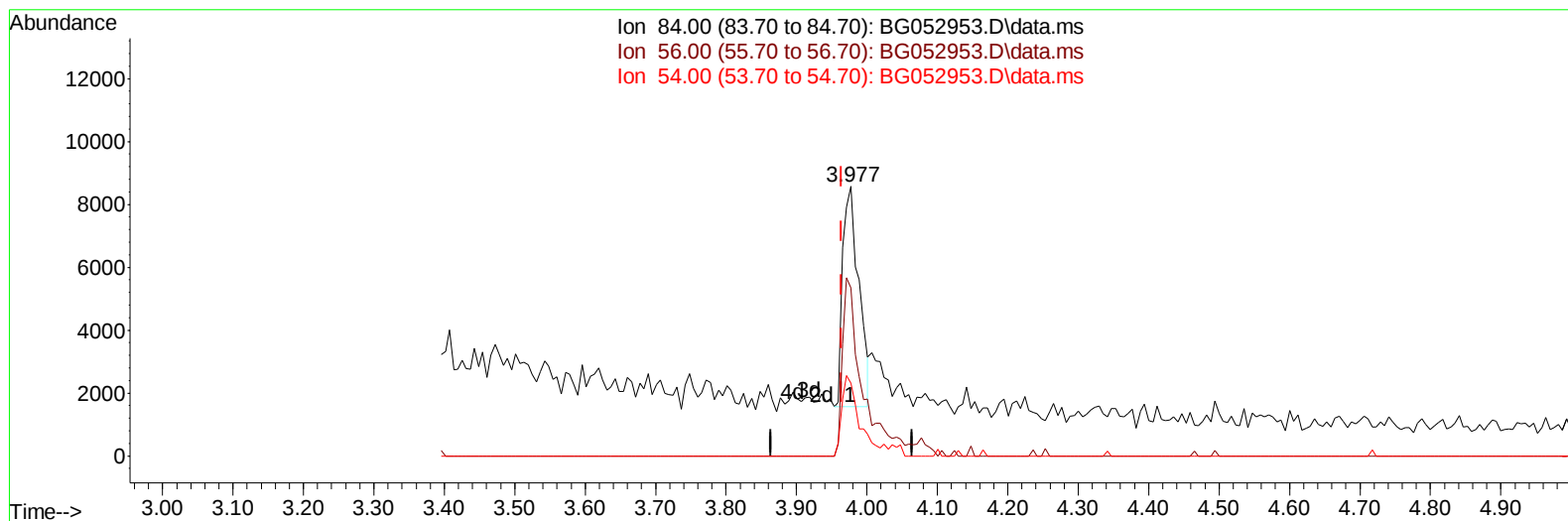
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG033022\
Data File : BG052953.D
Acq On : 30 Mar 2022 17:26
Operator : CG/JU
Sample : N2068-17
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
YAZ58

Manual IntegrationsAPPROVED

Quant Time: Mar 31 00:53:10 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG031822.M
Quant Title : SVOA CALIBRATION
QLast Update : Thu Mar 31 00:50:49 2022
Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 03/31/2022
Supervised By :Yogesh Patel 04/06/2022



TIC: BG052953.D\data.ms

(4) Pyridine-d5 (S)

3.977min (+ 0.013) 5.61 ng/u1

response 11006

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	65.20	62.34
54.00	28.70	27.03
0.00	0.00	0.00

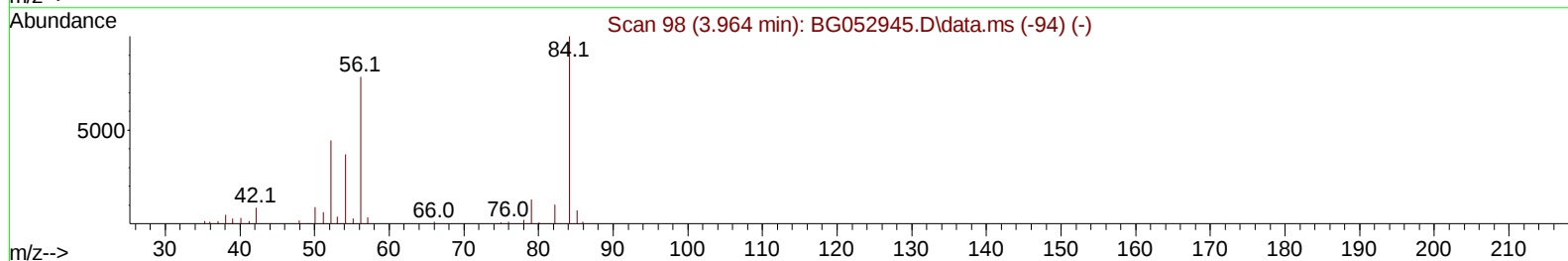
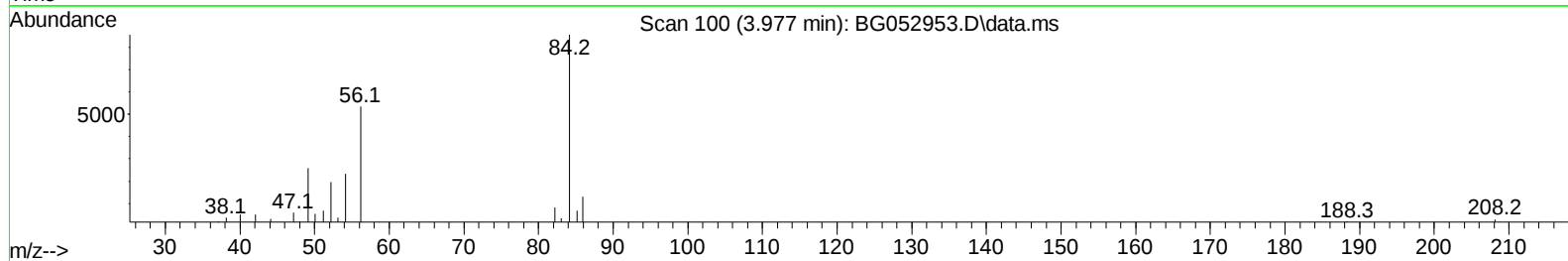
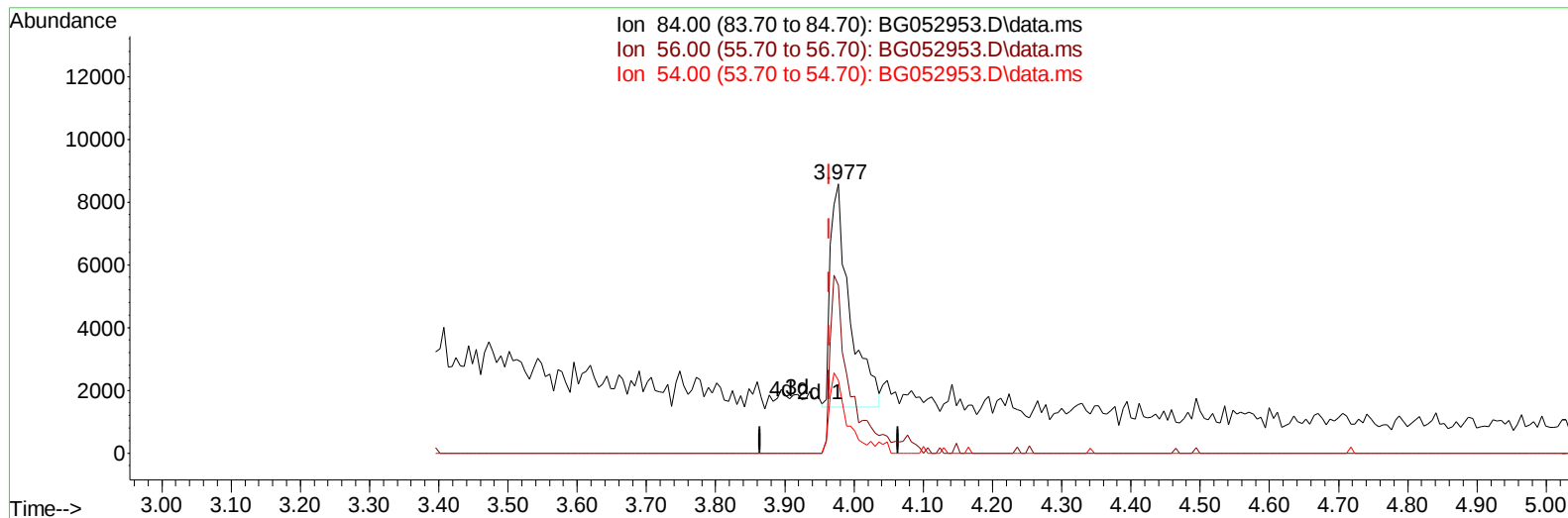
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TIC: BG052953.D\data.ms

(4) Pyridine-d5 (S)

3.977min (+ 0.013) 7.06 ng/u1 m

response 13863

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	65.20	62.34
54.00	28.70	27.03
0.00	0.00	0.00

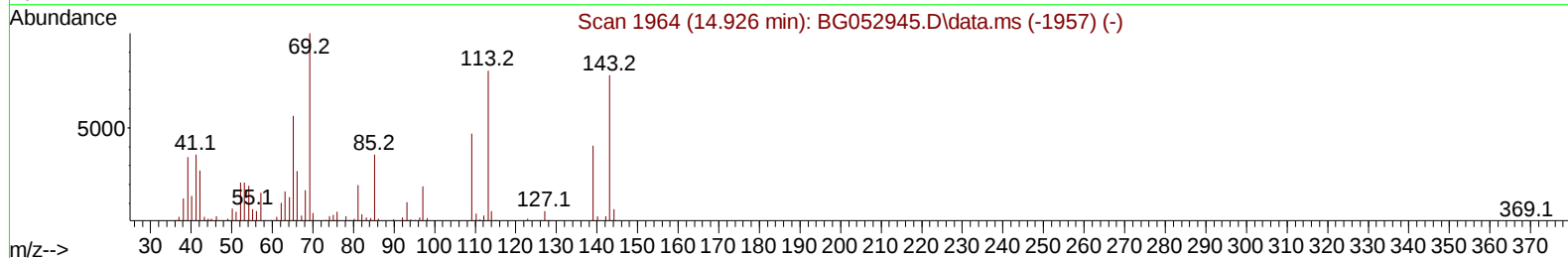
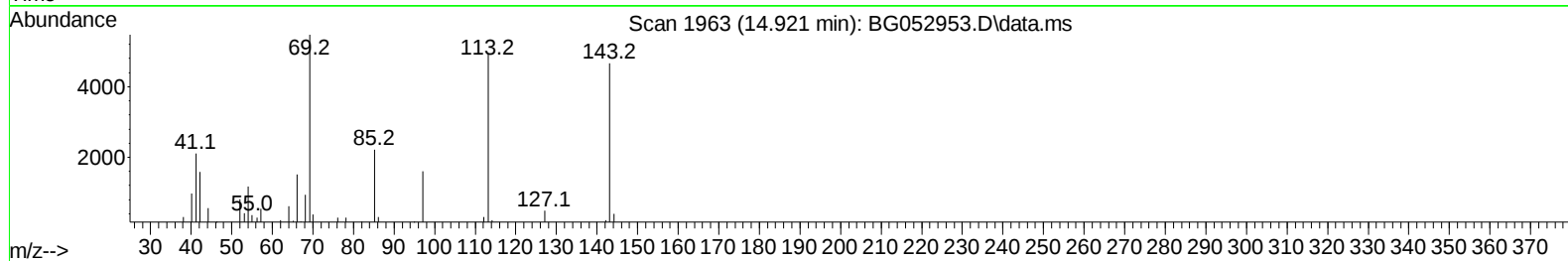
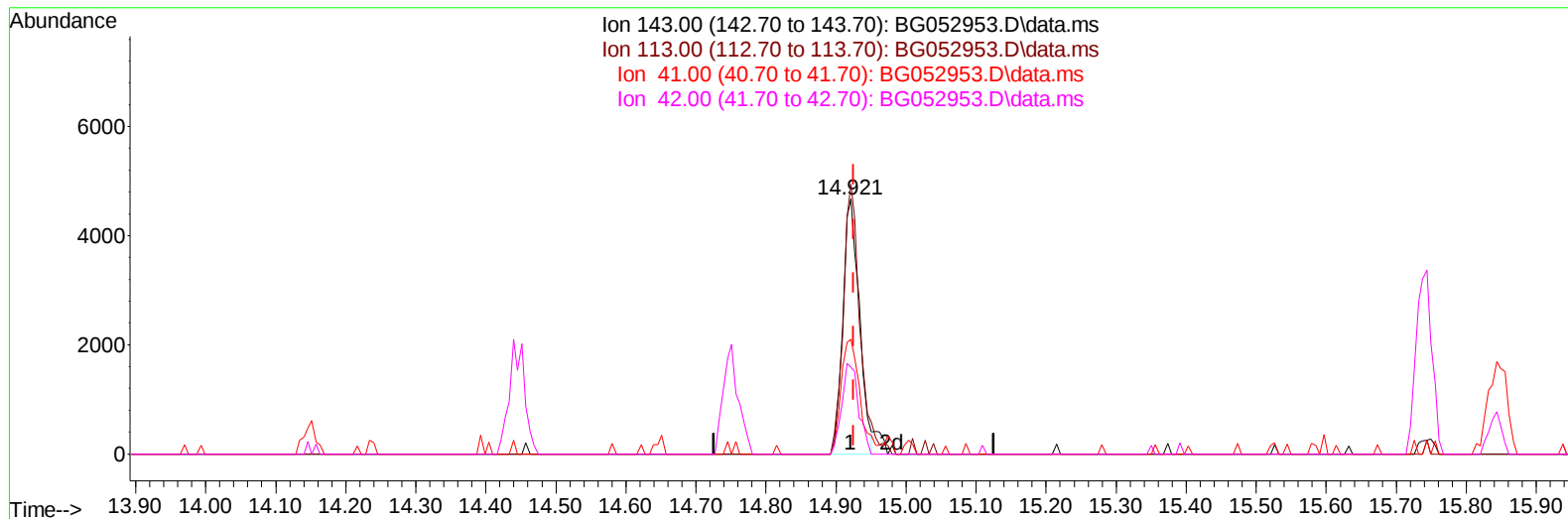
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Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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Manual IntegrationsAPPROVED

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TIC: BG052953.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.921min (-0.004) 7.37 ng/u1

response 7701

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	105.70	105.69
41.00	47.20	45.03
42.00	29.40	34.05

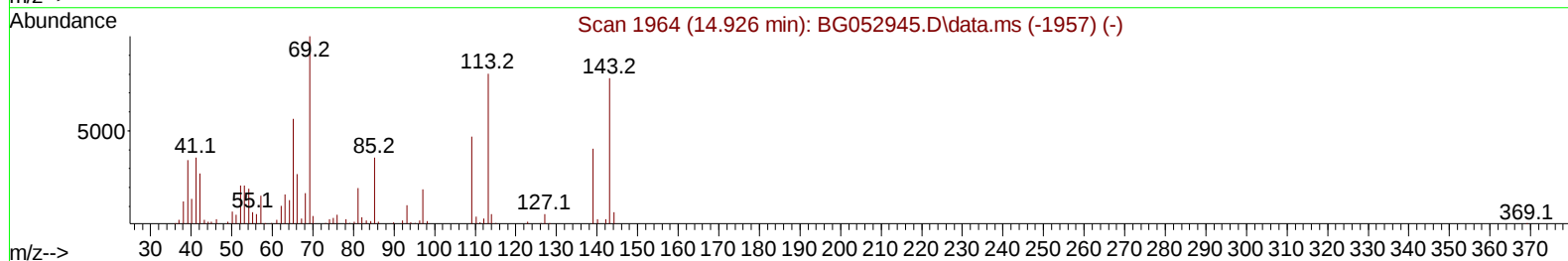
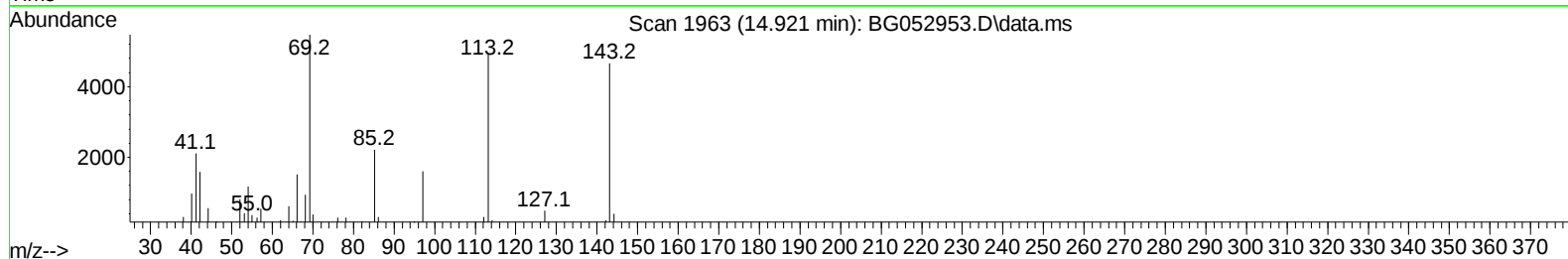
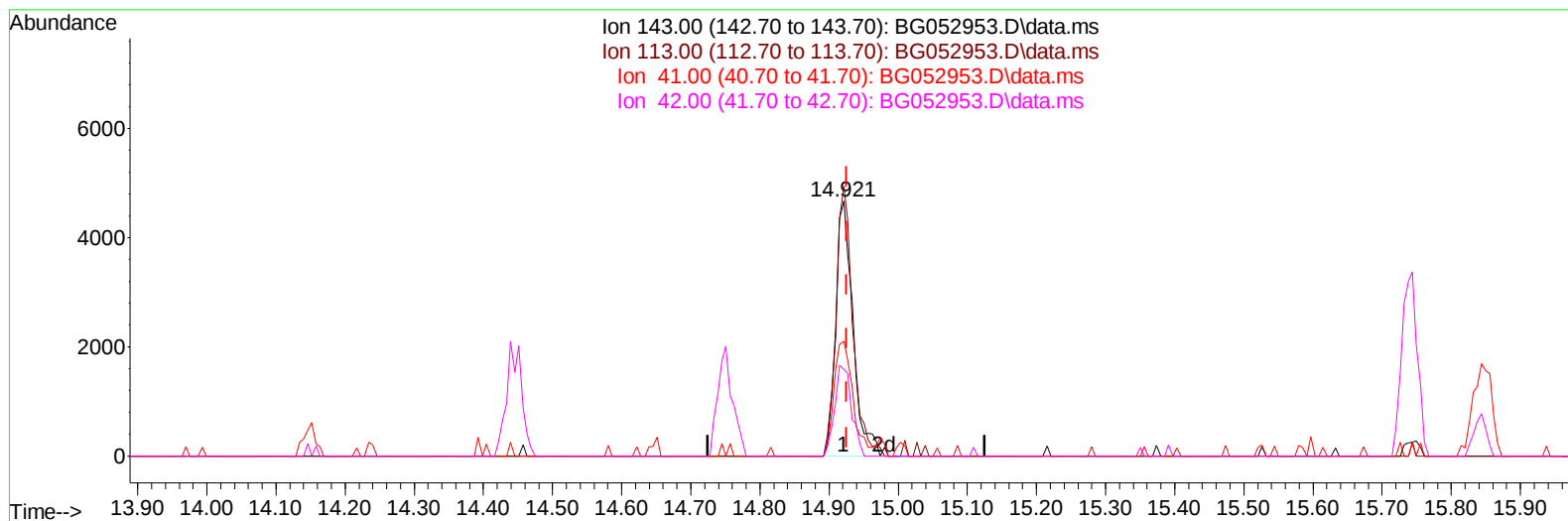
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG033022\
Data File : BG052953.D
Acq On : 30 Mar 2022 17:26
Operator : CG/JU
Sample : N2068-17
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
YAZ58

Manual IntegrationsAPPROVED

Quant Time: Mar 31 00:53:10 2022
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TIC: BG052953.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.921min (-0.004) 7.73 ng/ul m

response 8083

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	105.70	105.69
41.00	47.20	45.03
42.00	29.40	34.05

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.136	152	27764	20.000	ng/ul	0.00
20) Naphthalene-d8	10.944	136	110573	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.751	164	80925	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.494	188	192126	20.000	ng/ul	0.00
79) Chrysene-d12	21.777	240	229520	20.000	ng/ul	0.00
88) Perylene-d12	25.084	264	227118	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.554	96	3221	4.098	ng/uL	0.00
4) Pyridine-d5	3.977	84	13863m	7.065	ng/ul	0.01
7) Phenol-d5	7.279	99	19089	7.676	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.449	67	48145	31.700	ng/ul	0.00
11) 2-Chlorophenol-d4	7.660	132	39792	23.567	ng/ul	0.00
15) 4-Methylphenol-d8	8.829	113	31970	17.748	ng/ul	0.00
21) Nitrobenzene-d5	9.288	128	27820	34.952	ng/ul	0.00
24) 2-Nitrophenol-d4	10.016	143	28591	34.056	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.557	165	52463	28.725	ng/ul	0.00
31) 4-Chloroaniline-d4	11.068	131	56963	23.598	ng/ul	0.00
46) Dimethylphthalate-d6	14.146	166	212348	34.156	ng/ul	0.00
49) Acenaphthylene-d8	14.445	160	233139	30.861	ng/ul	0.00
54) 4-Nitrophenol-d4	14.921	143	8083m	7.733	ng/ul	0.00
60) Fluorene-d10	15.738	176	185007	33.891	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.844	200	35278	36.082	ng/ul	0.00
73) Anthracene-d10	17.594	188	319152	35.673	ng/ul	0.00
81) Pyrene-d10	19.873	212	422359	32.894	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.861	264	434515	37.027	ng/ul	0.00

Target Compounds Qvalue

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

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