

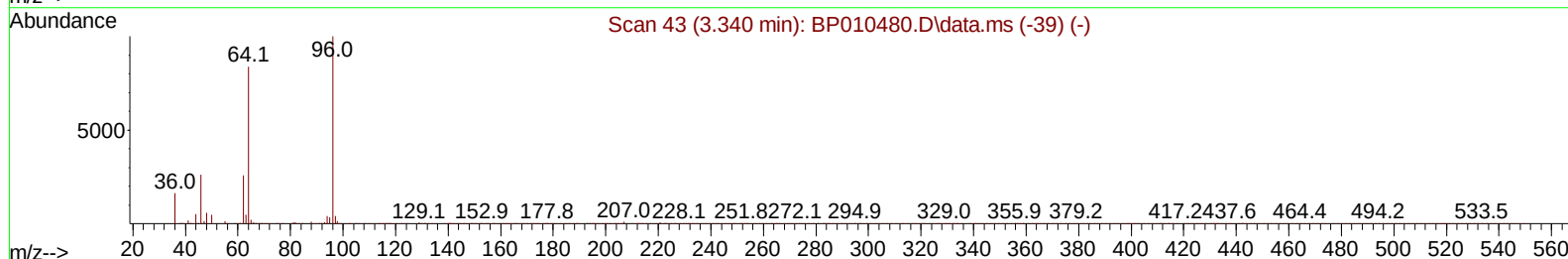
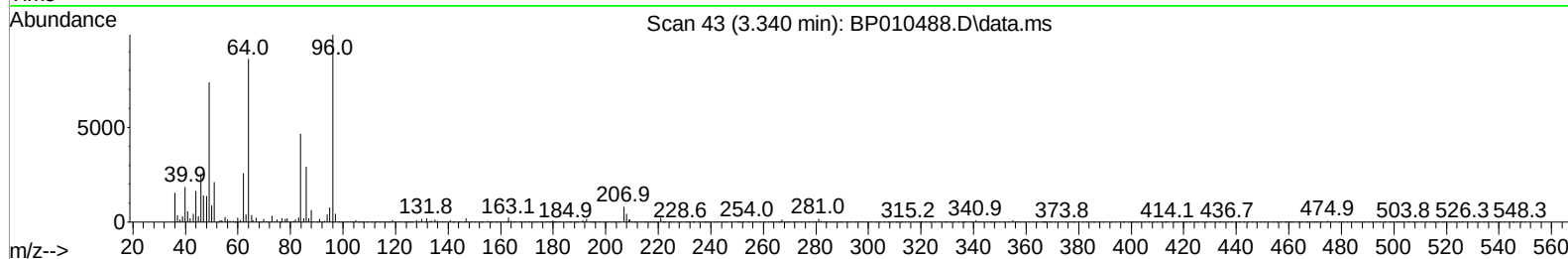
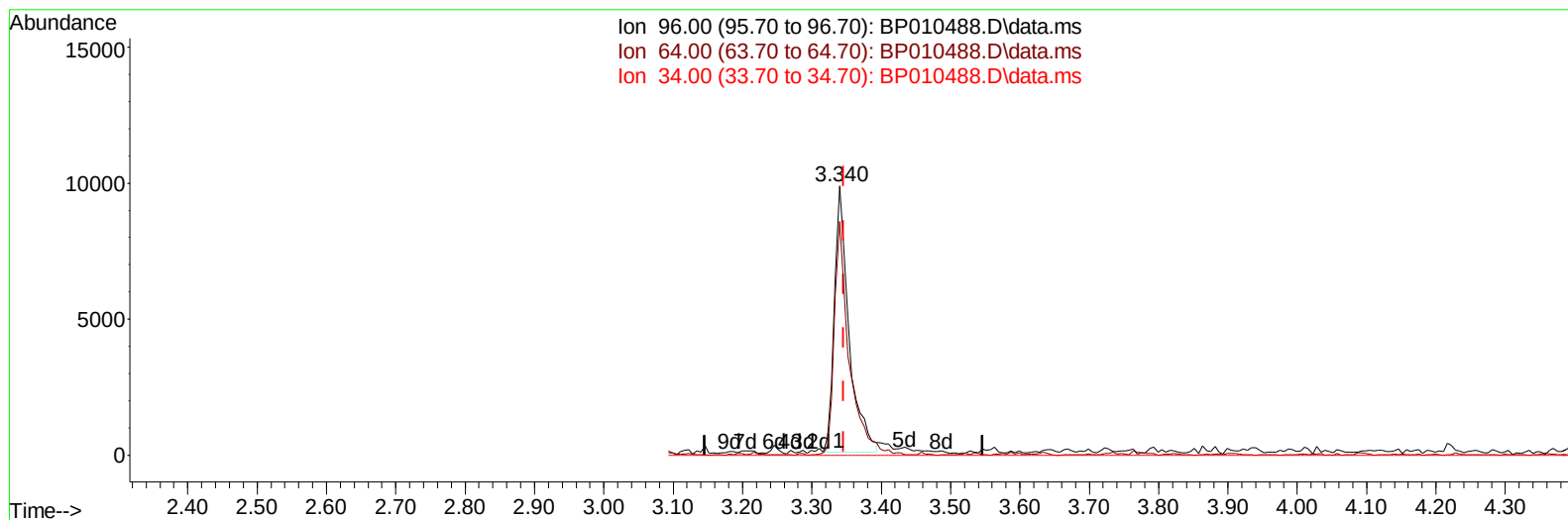
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP052322\
 Data File : BP010488.D
 Acq On : 23 May 2022 21:42
 Operator : CG/JU
 Sample : N2955-05
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0AA4

Manual IntegrationsAPPROVED

Quant Time: May 23 23:21:23 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP051322.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri May 13 15:04:31 2022
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 05/24/2022
 Supervised By :Yogesh Patel 05/26/2022



TIC: BP010488.D\data.ms

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

3.340min (-0.006) 4.00 ng/uL

response 14352

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	86.98#
34.00	0.00	0.00
0.00	0.00	0.00

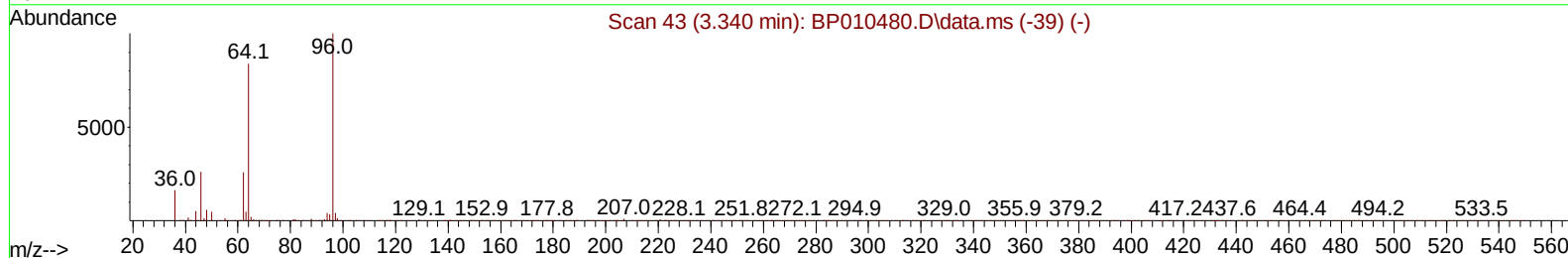
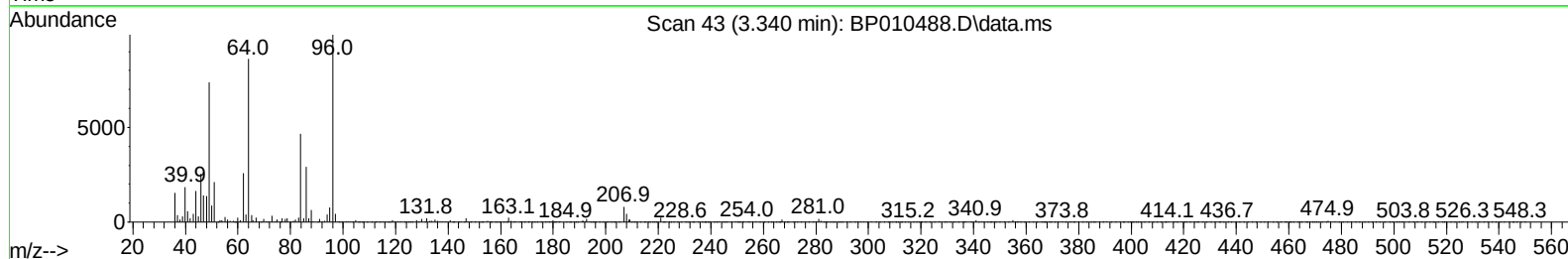
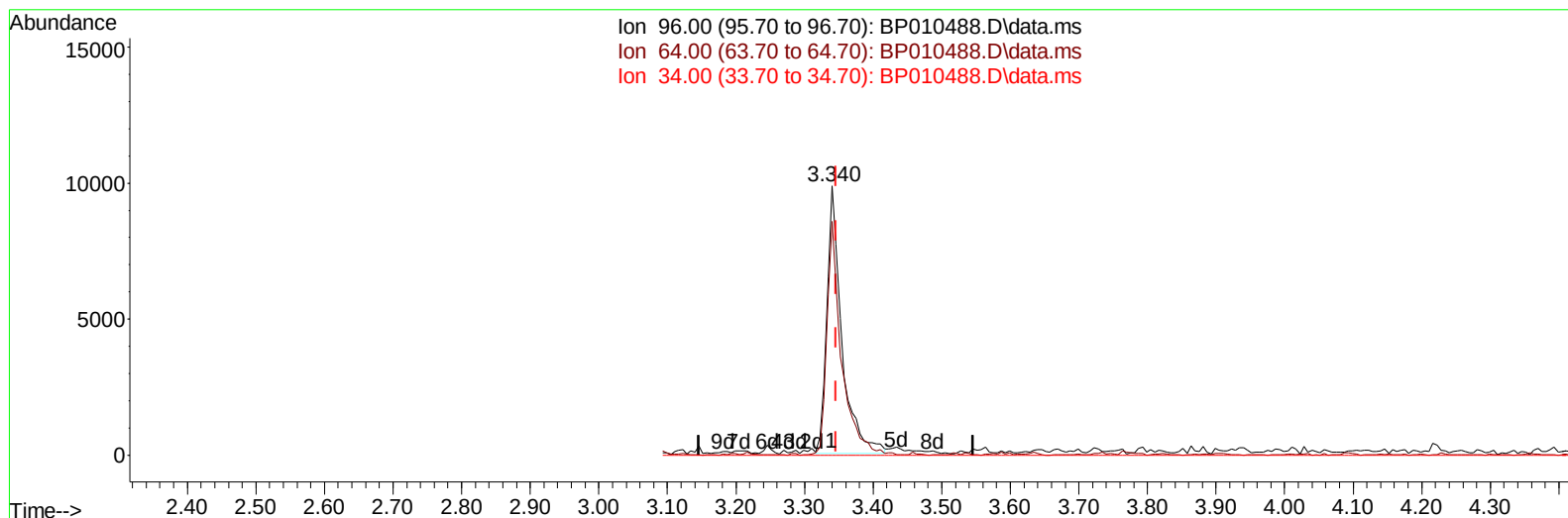
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TIC: BP010488.D\data.ms

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

3.340min (-0.006) 4.17 ng/uL m

response 14954

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	86.98#
34.00	0.00	0.00
0.00	0.00	0.00

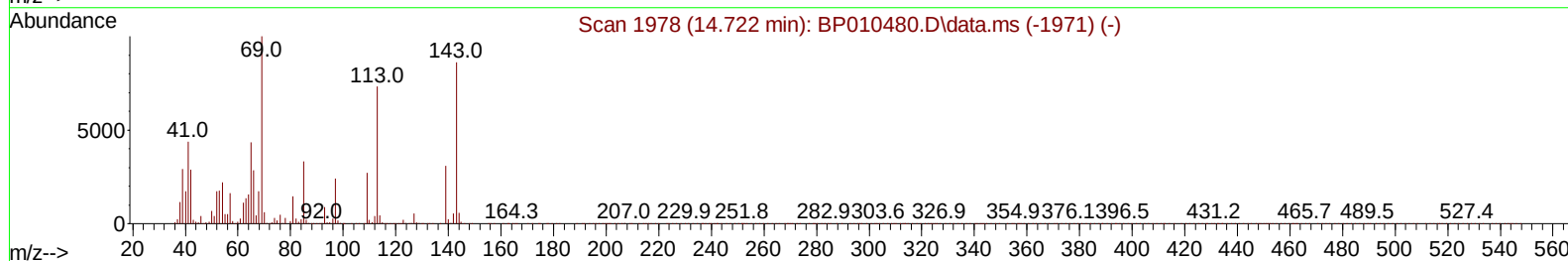
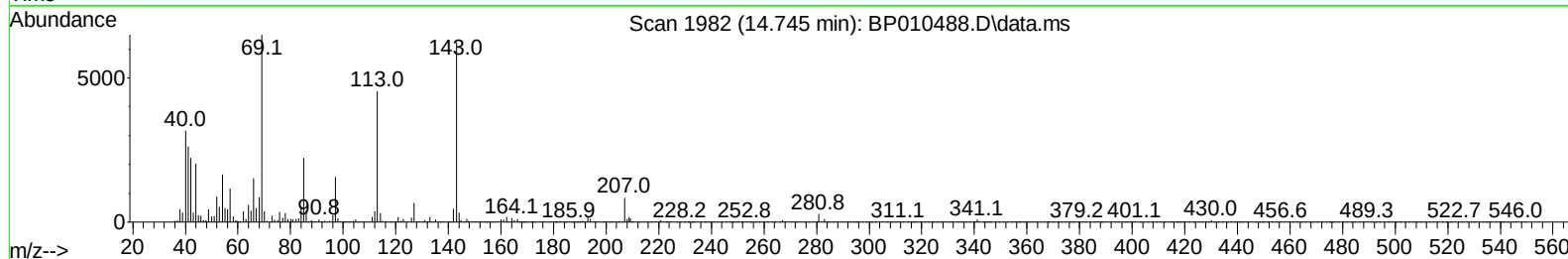
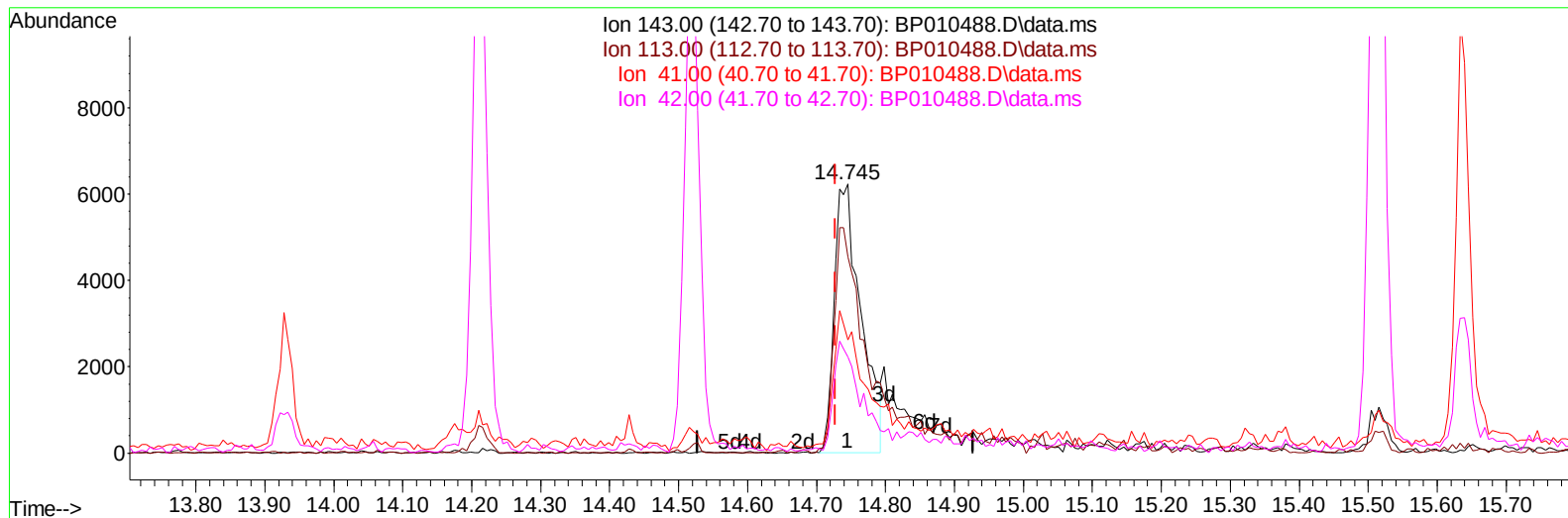
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ALS Vial : 17 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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TIC: BP010488.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.745min (+ 0.018) 2.75 ng/ul

response 16842

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	132.70	72.90#
41.00	23.20	42.00#
42.00	18.00	35.67#

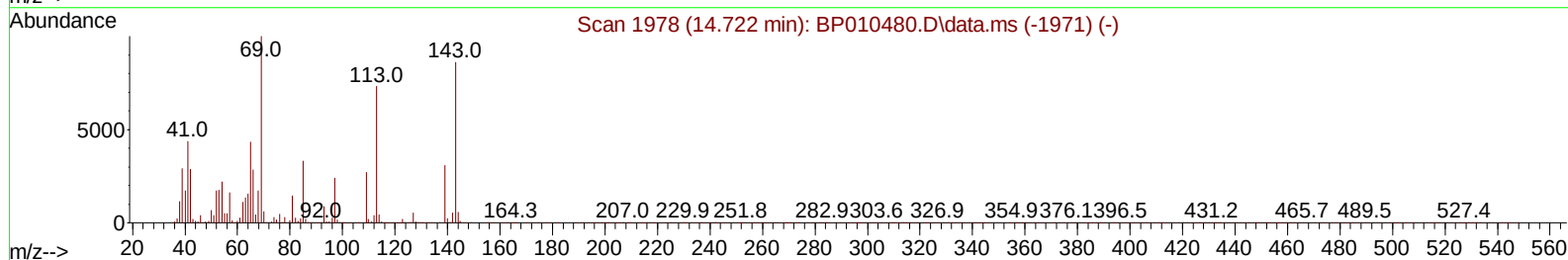
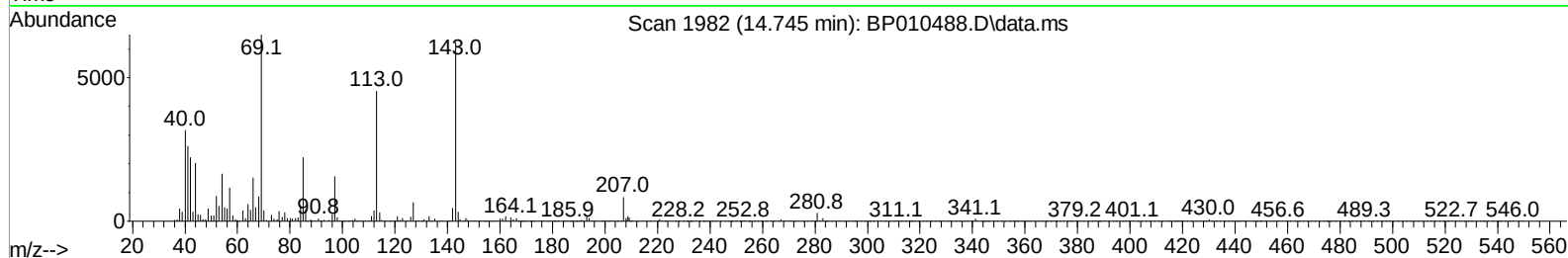
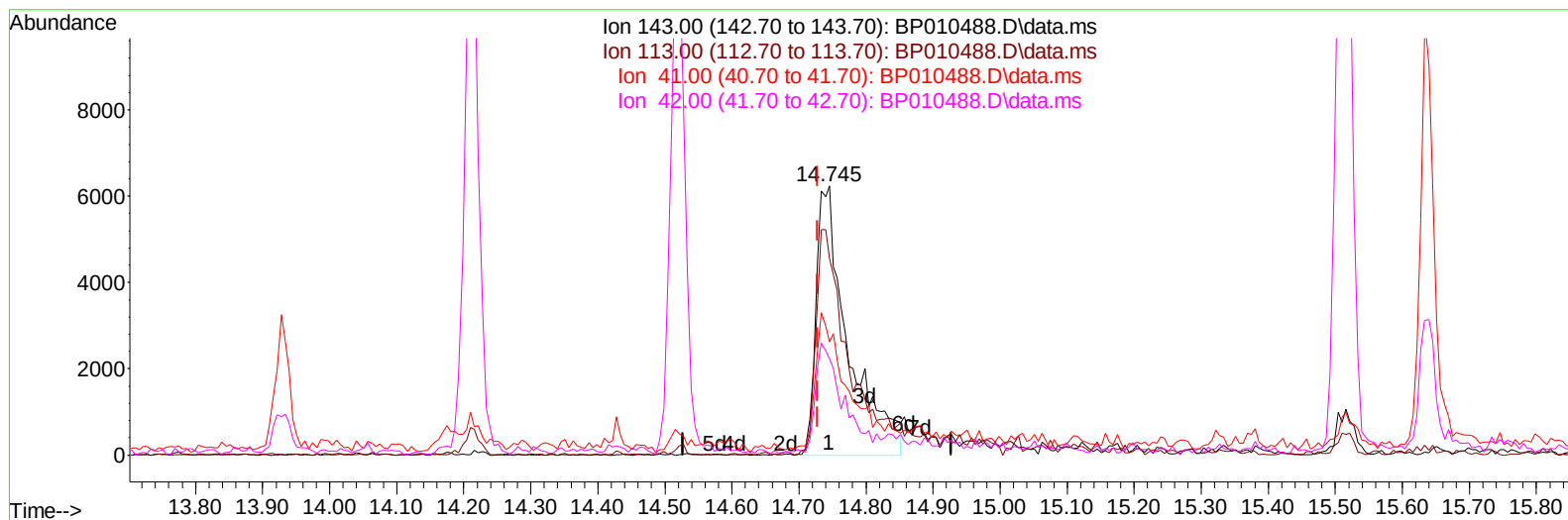
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 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
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 ClientSampleId :
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Manual IntegrationsAPPROVED

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TIC: BP010488.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.745min (+ 0.018) 3.38 ng/ul m

response 20732

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	132.70	72.90#
41.00	23.20	42.00#
42.00	18.00	35.67#

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.887	152	152981	20.000	ng/ul	-0.01
20) Naphthalene-d8	10.687	136	694850	20.000	ng/ul	-0.02
38) Acenaphthene-d10	14.522	164	467697	20.000	ng/ul	-0.01
64) Phenanthrene-d10	17.275	188	990077	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.357	240	845921	20.000	ng/ul	# 0.00
88) Perylene-d12	23.774	264	662603	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.340	96	14954m	4.167	ng/uL	0.00
4) Pyridine-d5	3.769	84	55300	5.427	ng/ul	0.00
7) Phenol-d5	7.057	99	72743	5.665	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.216	67	248754	28.835	ng/ul	-0.01
11) 2-Chlorophenol-d4	7.422	132	222826	22.679	ng/ul	-0.01
15) 4-Methylphenol-d8	8.599	113	148613	13.619	ng/ul	-0.01
21) Nitrobenzene-d5	9.051	128	157042	29.203	ng/ul	0.00
24) 2-Nitrophenol-d4	9.769	143	165373	29.018	ng/ul	-0.01
28) 2,4-Dichlorophenol-d3	10.310	165	266891	24.767	ng/ul	-0.01
31) 4-Chloroaniline-d4	10.828	131	343115	21.414	ng/ul	-0.01
46) Dimethylphthalate-d6	13.934	166	1033087	29.981	ng/ul	0.00
49) Acenaphthylene-d8	14.210	160	1197937	30.139	ng/ul	-0.01
54) 4-Nitrophenol-d4	14.745	143	20732m	3.380	ng/ul	0.02
60) Fluorene-d10	15.516	176	909916	30.353	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.639	200	136130	23.510	ng/ul	0.00
73) Anthracene-d10	17.369	188	1565044	34.909	ng/ul	-0.01
81) Pyrene-d10	19.604	212	1796282	39.985	ng/ul	-0.01
92) Benzo(a)pyrene-d12	23.615	264	1175054	35.317	ng/ul	-0.02
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.375	88	10726	2.885	ng/ul#	21

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

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