

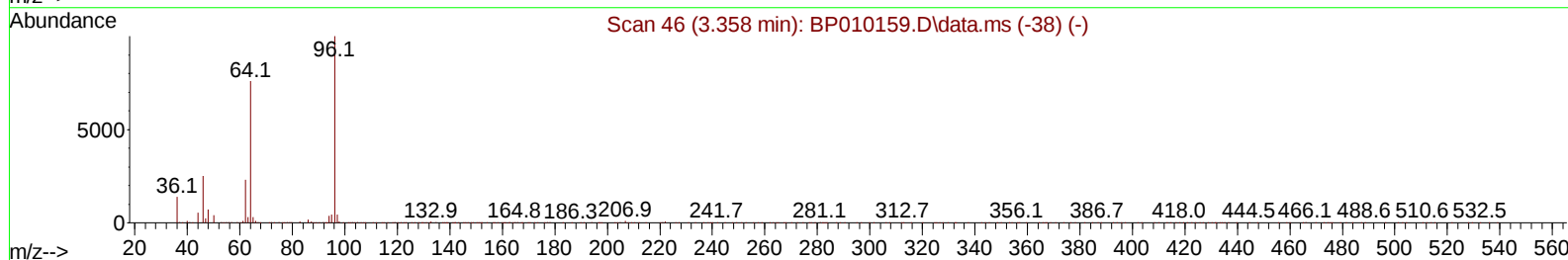
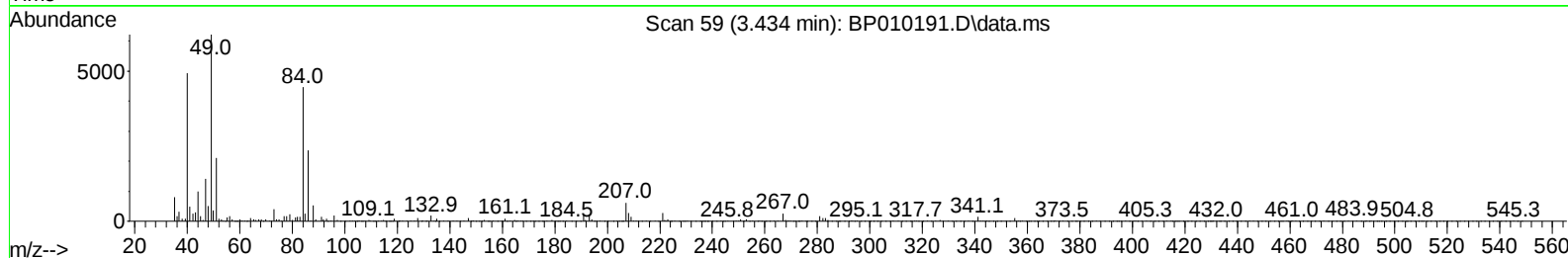
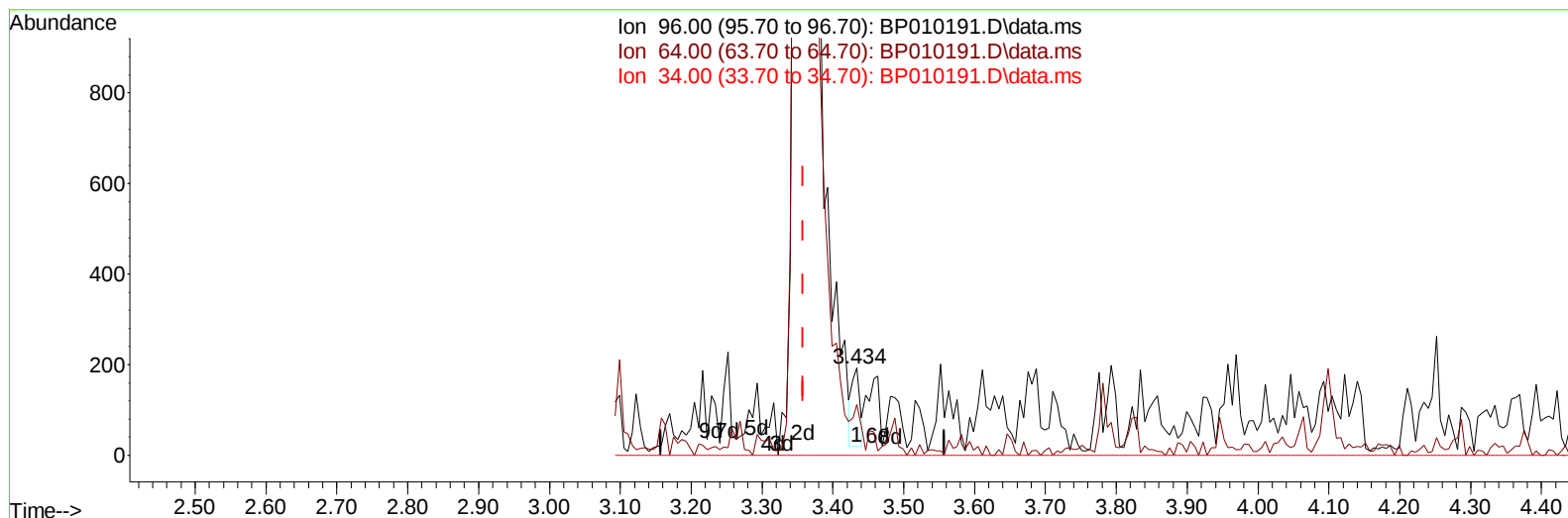
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP050222\
 Data File : BP010191.D
 Acq On : 03 May 2022 01:57
 Operator : CG/JU
 Sample : N2615-05
 Misc :
 ALS Vial : 27 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BH0D2

Manual IntegrationsAPPROVED

Quant Time: May 03 02:54:12 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: BP010191.D\data.ms

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

3.434min (+ 0.076) 0.05 ng/uL

response 136

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	58.03
34.00	0.00	0.00
0.00	0.00	0.00

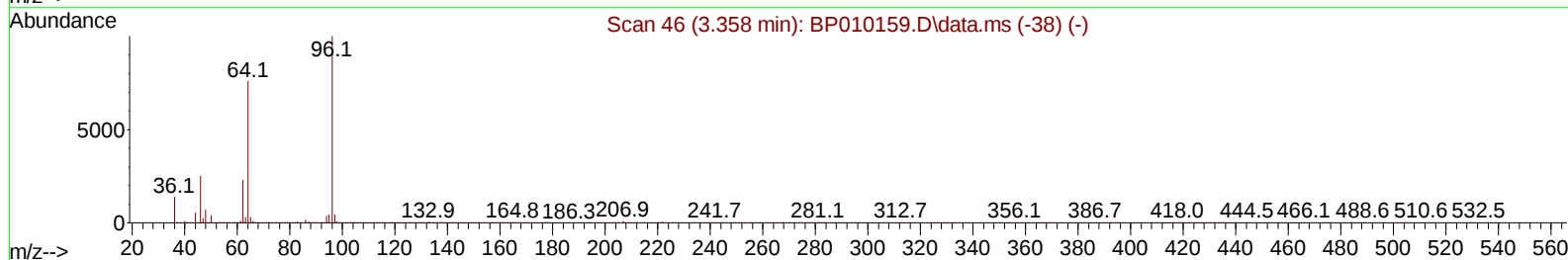
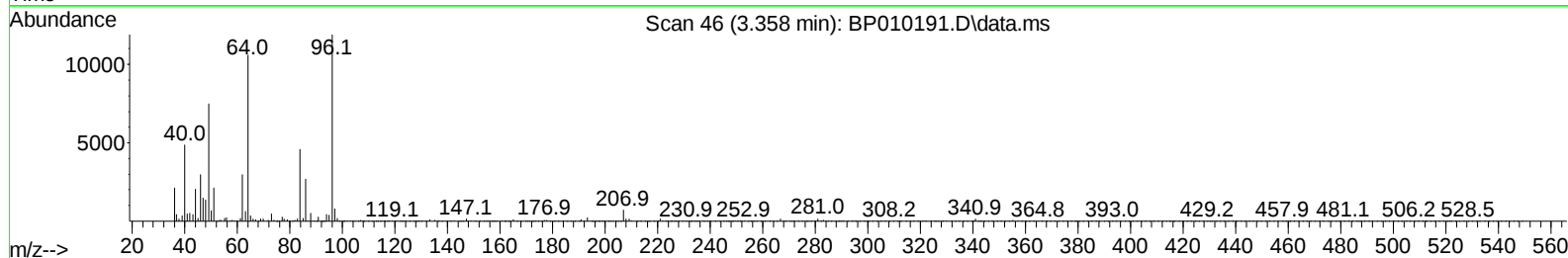
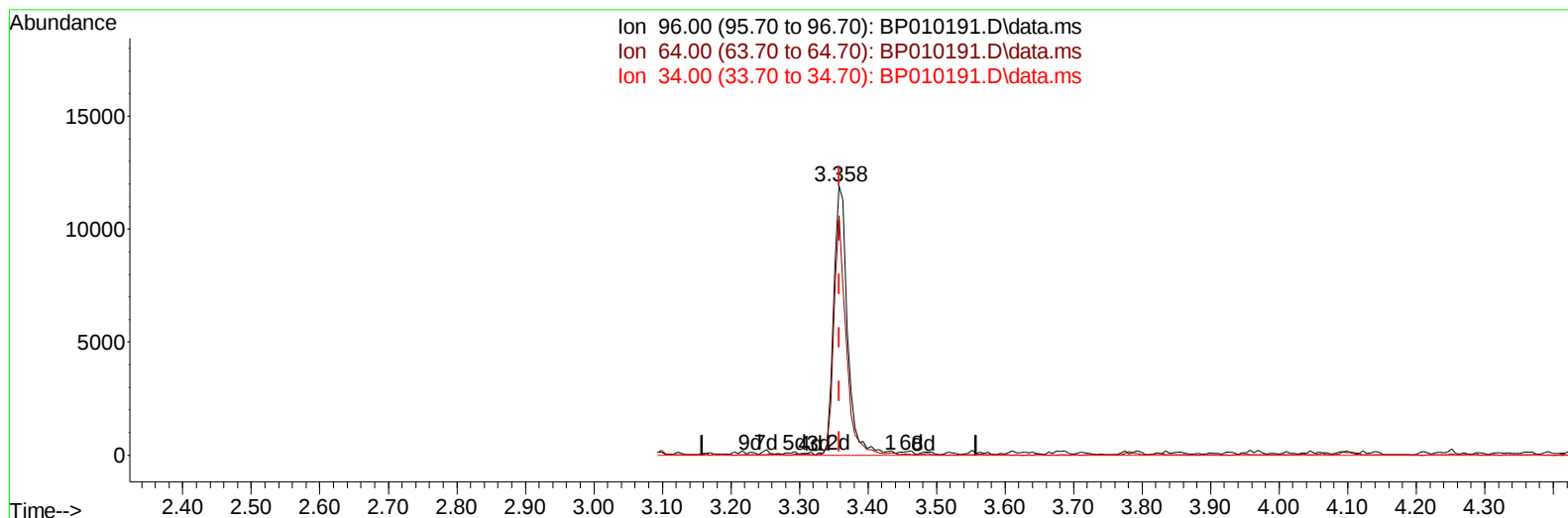
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TIC: BP010191.D\data.ms

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

3.358min (-0.000) 5.59 ng/uL m

response 16778

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	89.11#
34.00	0.00	0.00
0.00	0.00	0.00

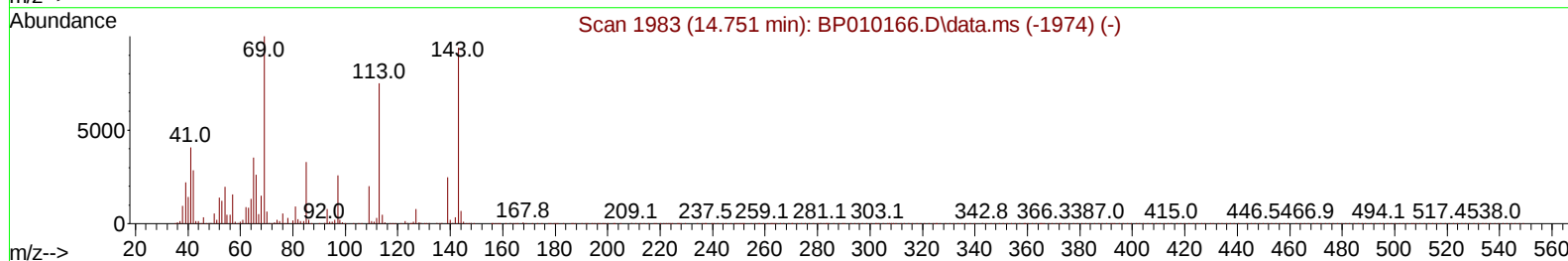
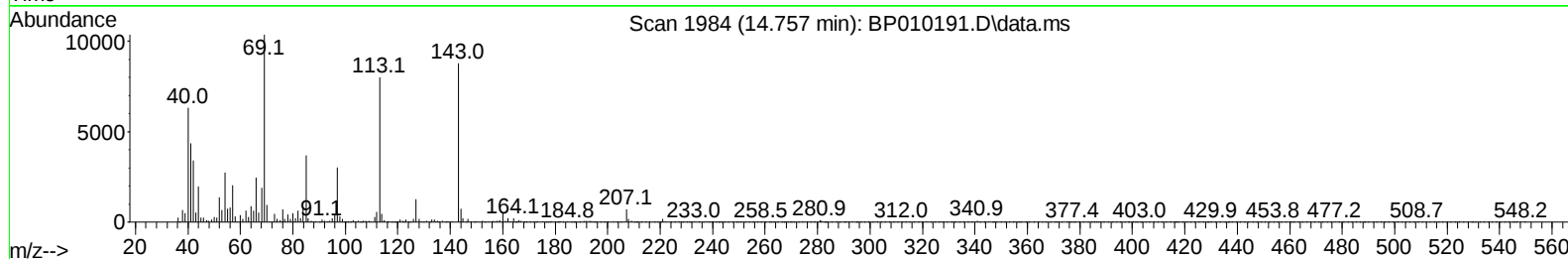
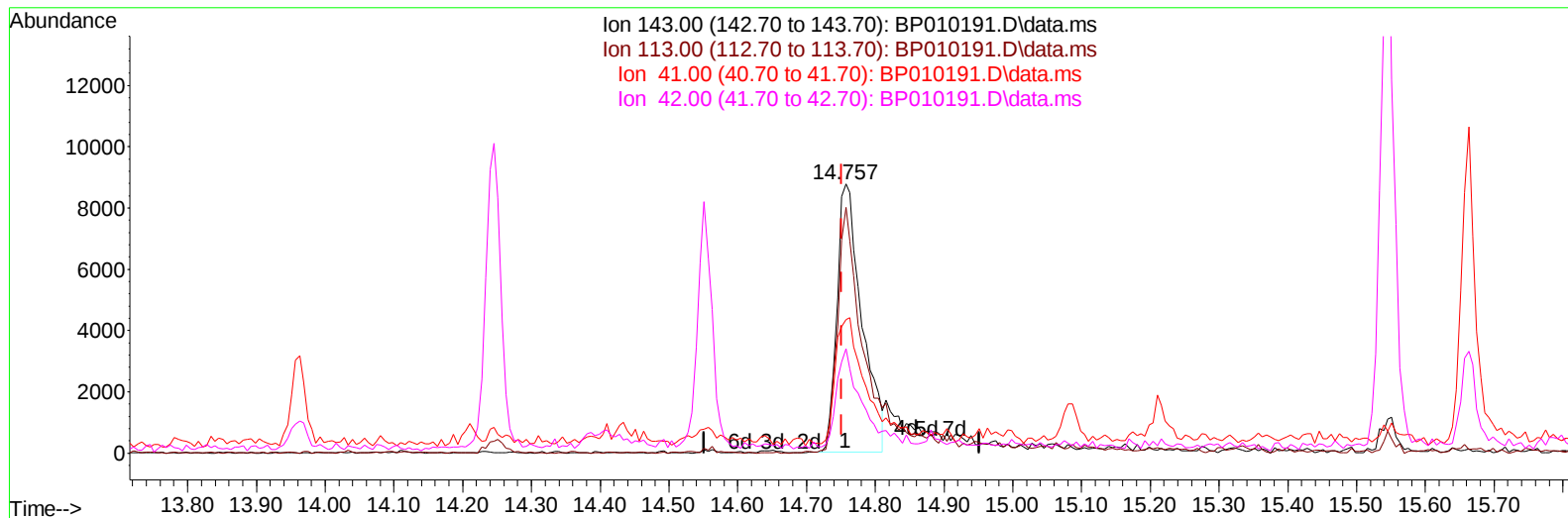
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Misc :
ALS Vial : 27 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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TIC: BP010191.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.757min (+ 0.006) 4.54 ng/ul

response 22103

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	132.70	91.33#
41.00	23.20	49.56#
42.00	18.00	38.72#

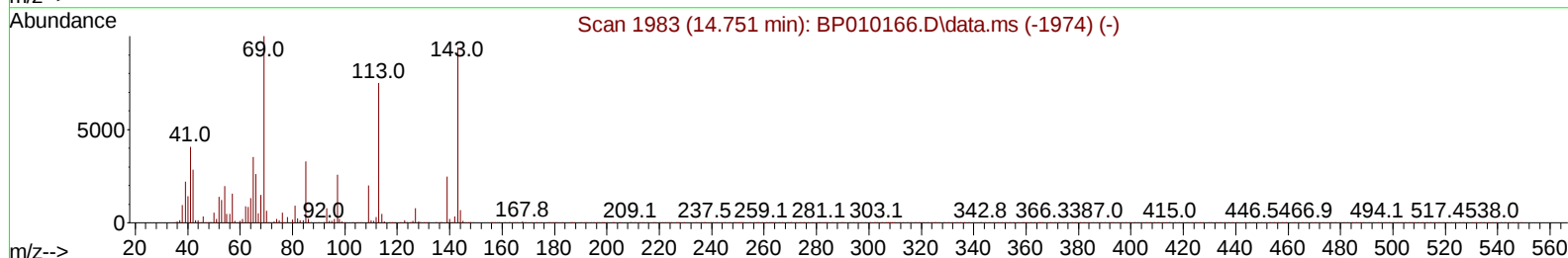
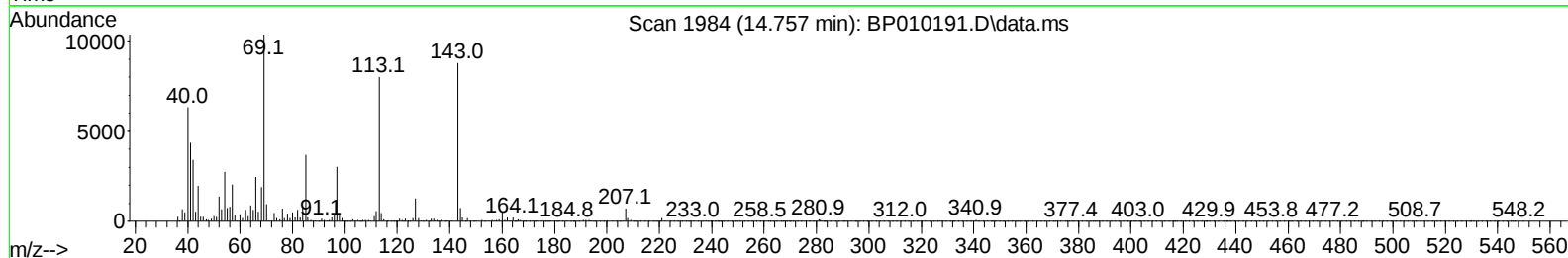
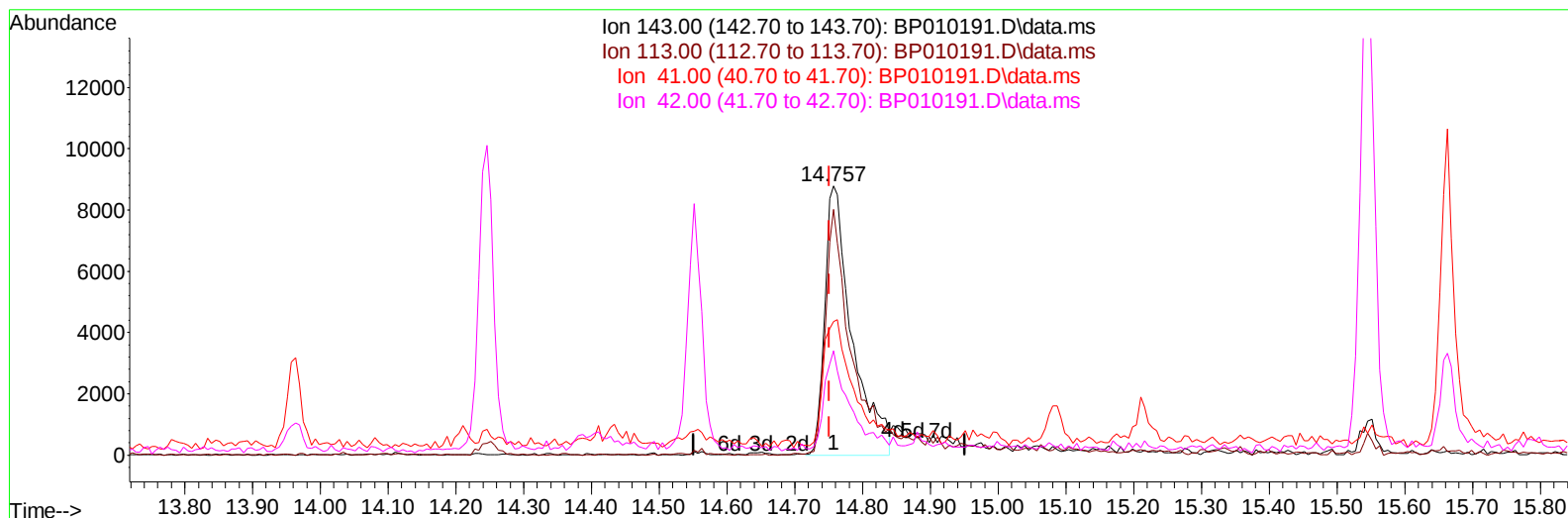
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 Operator : CG/JU
 Sample : N2615-05
 Misc :
 ALS Vial : 27 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
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Manual IntegrationsAPPROVED

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TIC: BP010191.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.757min (+ 0.006) 5.02 ng/ul m

response 24441

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	132.70	91.33#
41.00	23.20	49.56#
42.00	18.00	38.72#

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.922	152	130904	20.000	ng/ul	0.00
20) Naphthalene-d8	10.722	136	558895	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.551	164	340307	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.304	188	625693	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.386	240	516520	20.000	ng/ul	# 0.00
88) Perylene-d12	23.827	264	513695	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.358	96	16778m	5.589	ng/uL	0.00
4) Pyridine-d5	3.787	84	43385	5.137	ng/ul	0.01
7) Phenol-d5	7.087	99	83325	6.998	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.246	67	275831	38.607	ng/ul	0.00
11) 2-Chlorophenol-d4	7.452	132	249002	28.237	ng/ul	0.00
15) 4-Methylphenol-d8	8.628	113	171178	17.352	ng/ul	0.00
21) Nitrobenzene-d5	9.081	128	166805	37.525	ng/ul	0.00
24) 2-Nitrophenol-d4	9.804	143	177719	36.608	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.346	165	282554	29.897	ng/ul	0.00
31) 4-Chloroaniline-d4	10.857	131	327436	24.163	ng/ul	0.00
46) Dimethylphthalate-d6	13.963	166	1104283	41.715	ng/ul	0.00
49) Acenaphthylene-d8	14.245	160	1203424	37.562	ng/ul	0.00
54) 4-Nitrophenol-d4	14.757	143	24441m	5.015	ng/ul	0.00
60) Fluorene-d10	15.545	176	900312	39.537	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.663	200	148828	37.618	ng/ul	0.00
73) Anthracene-d10	17.404	188	1459504	48.216	ng/ul	0.00
81) Pyrene-d10	19.633	212	1598320	54.411	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.674	264	1387321	51.514	ng/ul	0.00

Target Compounds Qvalue

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

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