

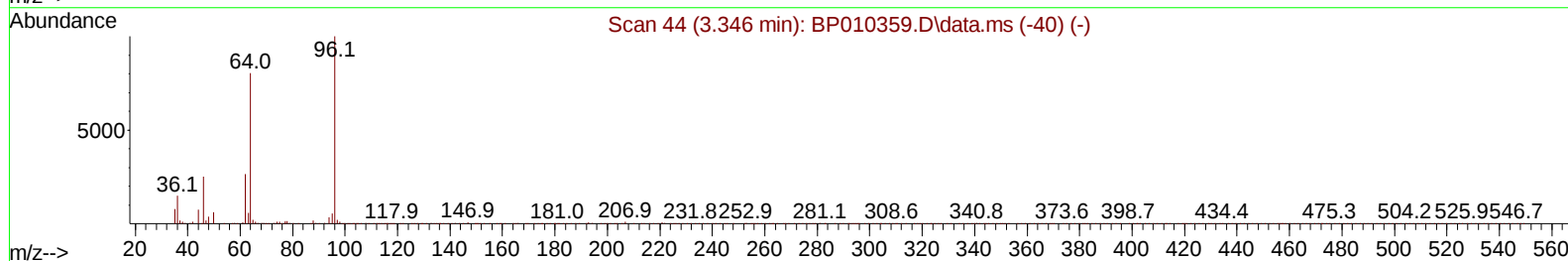
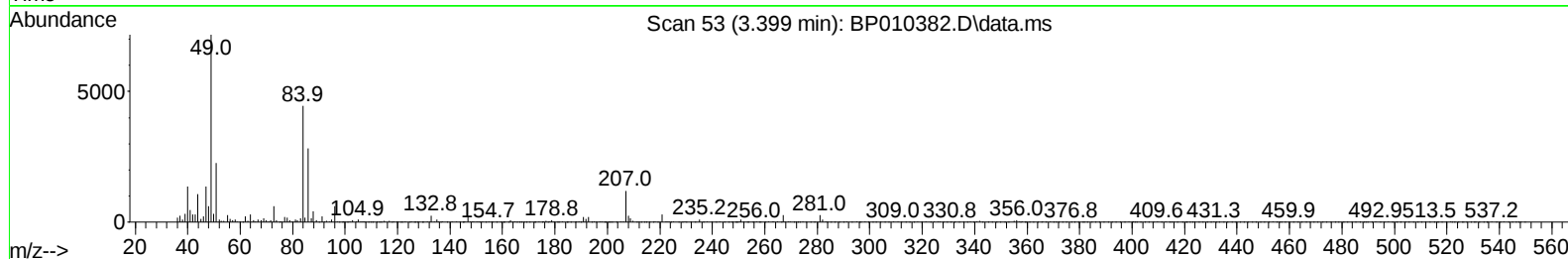
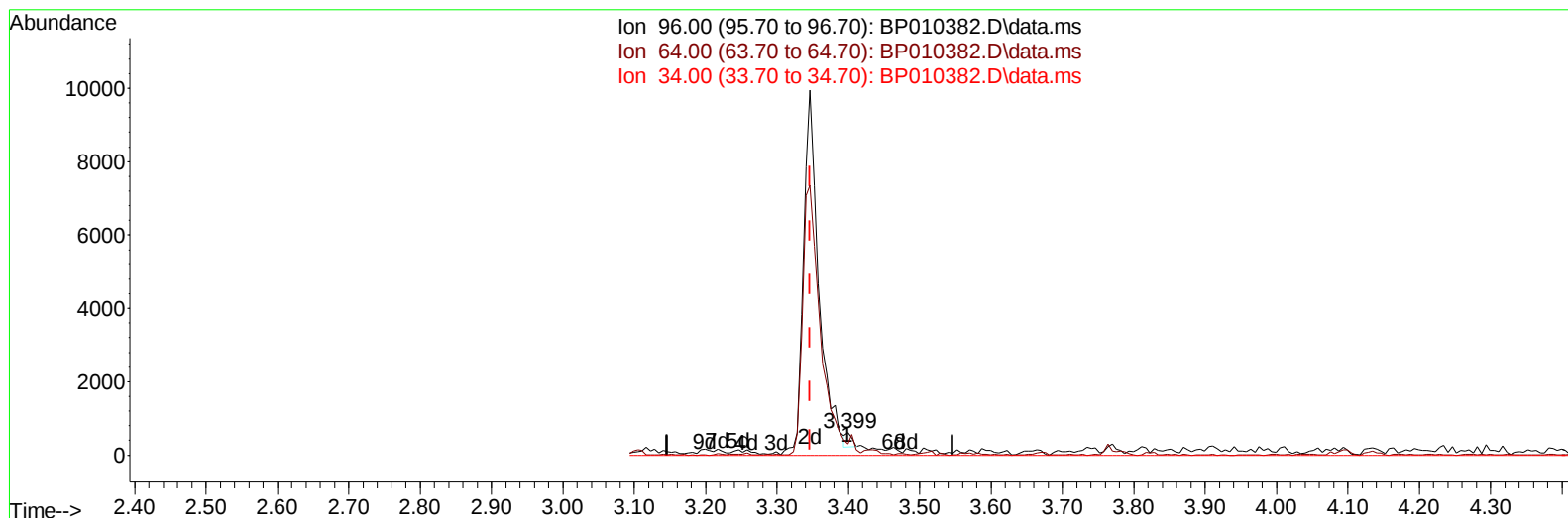
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051622\
 Data File : BP010382.D
 Acq On : 17 May 2022 05:56
 Operator : CG/JU
 Sample : N2831-08
 Misc :
 ALS Vial : 30 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 YB2L7

Manual IntegrationsAPPROVED

Quant Time: May 17 07:08:15 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 05/17/2022
 Supervised By :mohammad ahmed 05/18/2022



TIC: BP010382.D\data.ms

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

3.399min (+ 0.053) 0.06 ng/uL

response 225

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	48.37
34.00	0.00	0.00
0.00	0.00	0.00

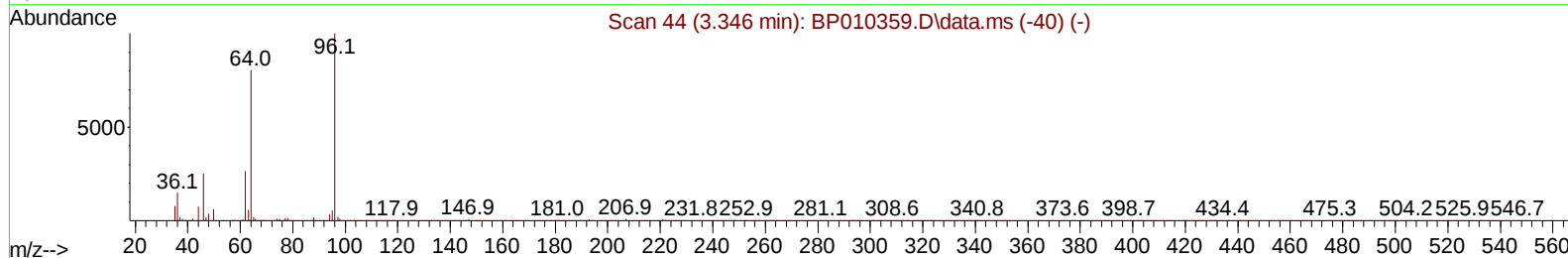
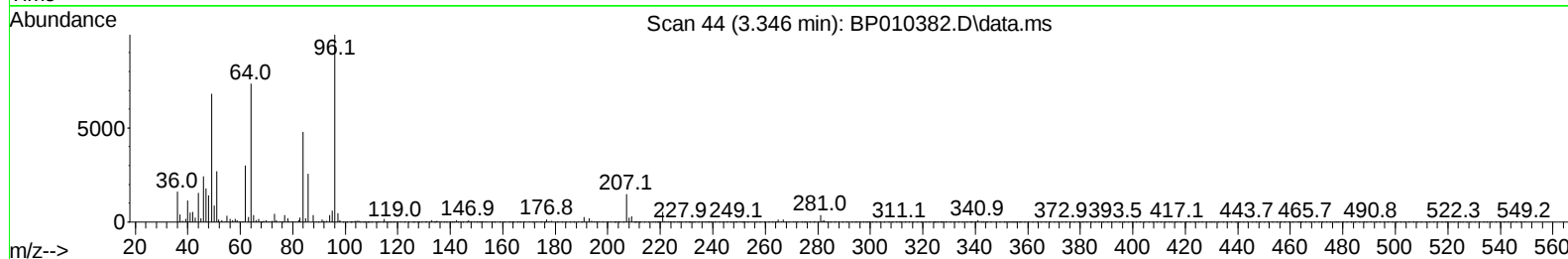
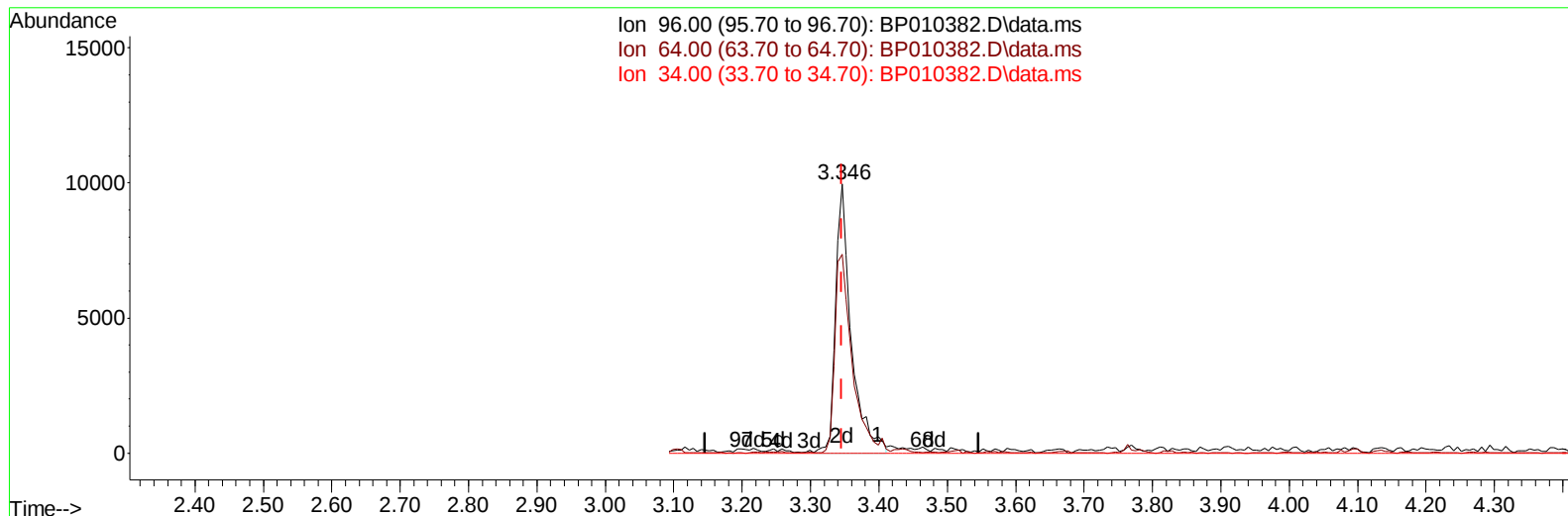
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(3) 1,4-Dioxane-d8 (S)

3.346min (+ 0.000) 4.06 ng/uL m

response 15874

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	73.83#
34.00	0.00	0.00
0.00	0.00	0.00

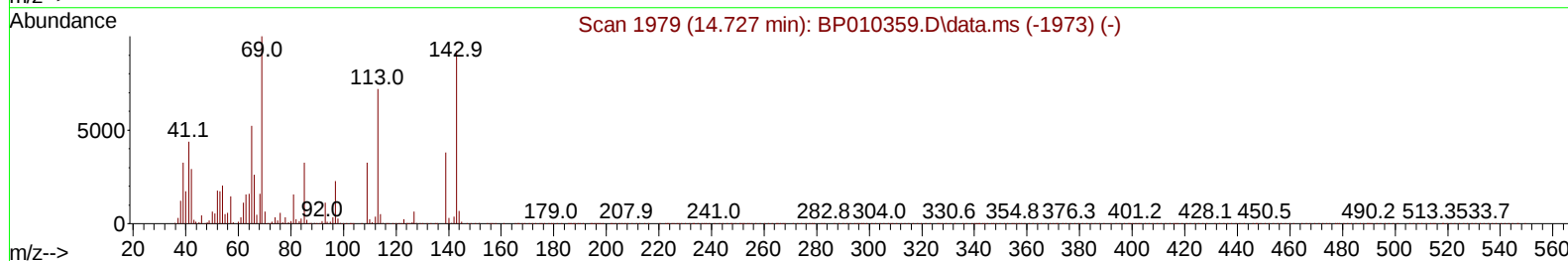
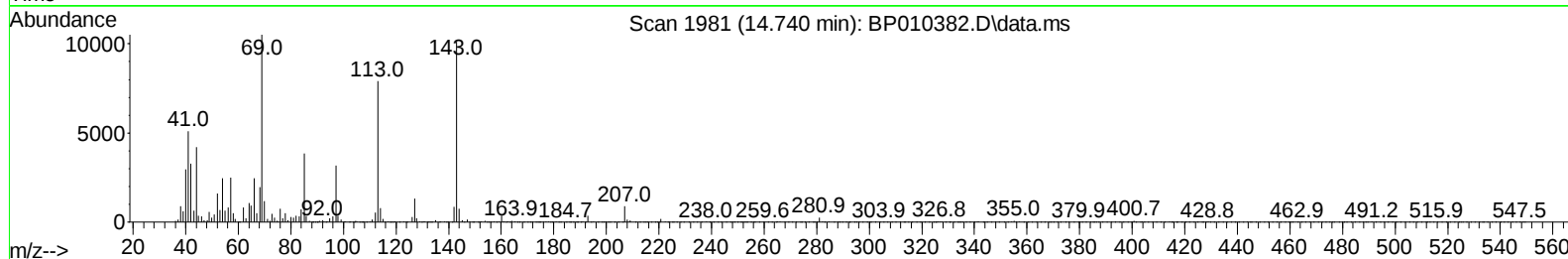
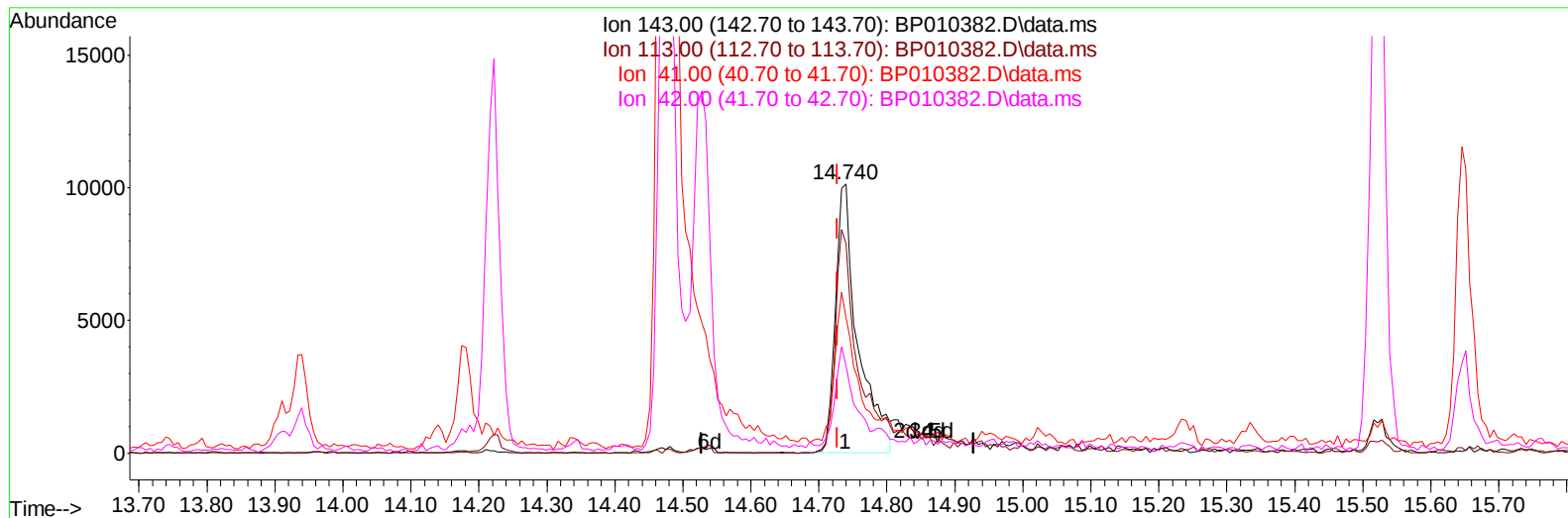
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TIC: BP010382.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.740min (+ 0.012) 3.69 ng/u1

response 23369

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	132.70	77.99#
41.00	23.20	50.36#
42.00	18.00	32.45#

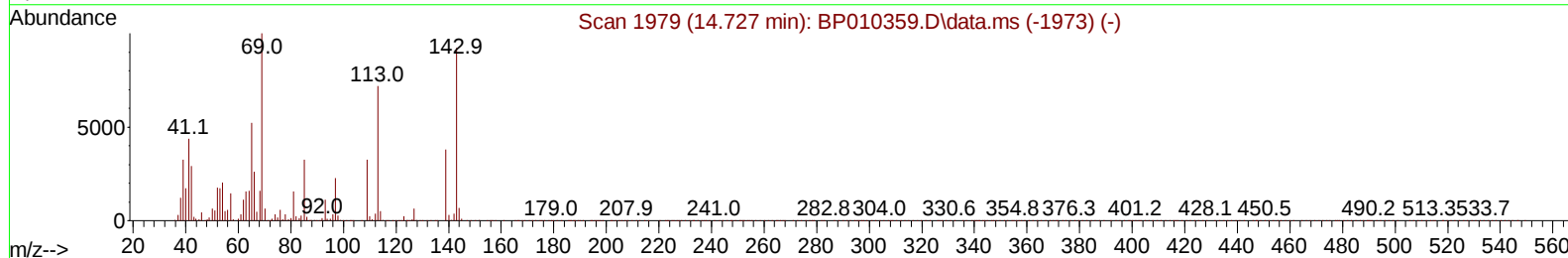
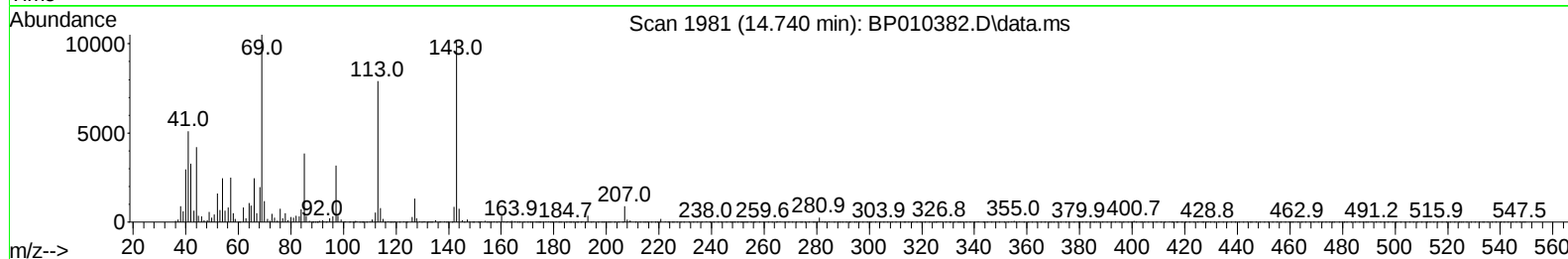
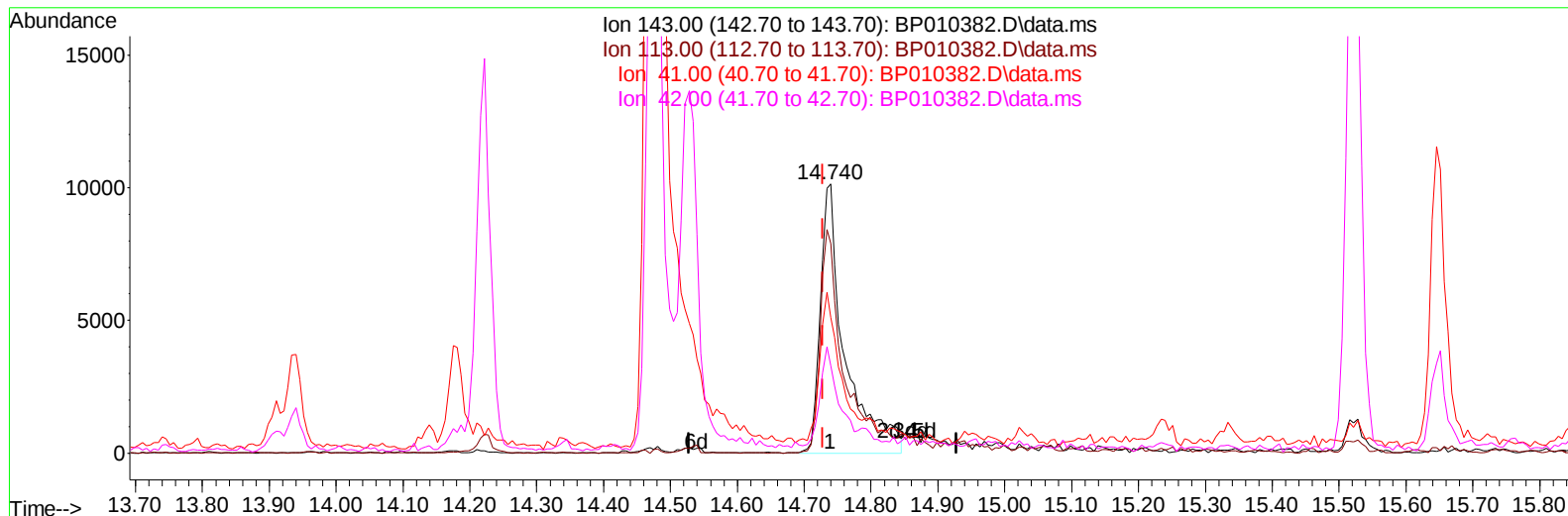
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Instrument :
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TIC: BP010382.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.740min (+ 0.012) 4.08 ng/ul m

response 25830

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	132.70	77.99#
41.00	23.20	50.36#
42.00	18.00	32.45#

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Compound	R.T.	Q	Ion	Response	Conc	Units	Dev(Min)

Internal Standards							
1) 1,4-Dichlorobenzene-d4	7.899	152		166691	20.000	ng/ul	0.00
20) Naphthalene-d8	10.699	136		724869	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.528	164		483122	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.281	188		1079483	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.363	240		1071883	20.000	ng/ul	# 0.00
88) Perylene-d12	23.792	264		966282	20.000	ng/ul	0.00
System Monitoring Compounds							
3) 1,4-Dioxane-d8	3.346	96		15874m	4.059	ng/uL	0.00
4) Pyridine-d5	3.770	84		95434	8.595	ng/ul	0.00
7) Phenol-d5	7.064	99		91353	6.530	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.222	67		306731	32.631	ng/ul	0.00
11) 2-Chlorophenol-d4	7.428	132		272089	25.416	ng/ul	0.00
15) 4-Methylphenol-d8	8.605	113		192732	16.210	ng/ul	0.00
21) Nitrobenzene-d5	9.058	128		188450	33.592	ng/ul	0.00
24) 2-Nitrophenol-d4	9.781	143		226592	38.113	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.322	165		349487	31.089	ng/ul	0.00
31) 4-Chloroaniline-d4	10.840	131		332954	19.920	ng/ul	0.00
46) Dimethylphthalate-d6	13.940	166		1280915	35.986	ng/ul	0.00
49) Acenaphthylene-d8	14.222	160		1446635	35.234	ng/ul	0.00
54) 4-Nitrophenol-d4	14.740	143		25830m	4.077	ng/ul	0.01
60) Fluorene-d10	15.522	176		1102566	35.605	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.651	200		164242	26.016	ng/ul	0.00
73) Anthracene-d10	17.381	188		1746864	35.737	ng/ul	0.00
81) Pyrene-d10	19.616	212		2138474	37.567	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.639	264		1736451	35.788	ng/ul	0.00

Target Compounds Qvalue

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

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