Quantitation Report (Qedit)

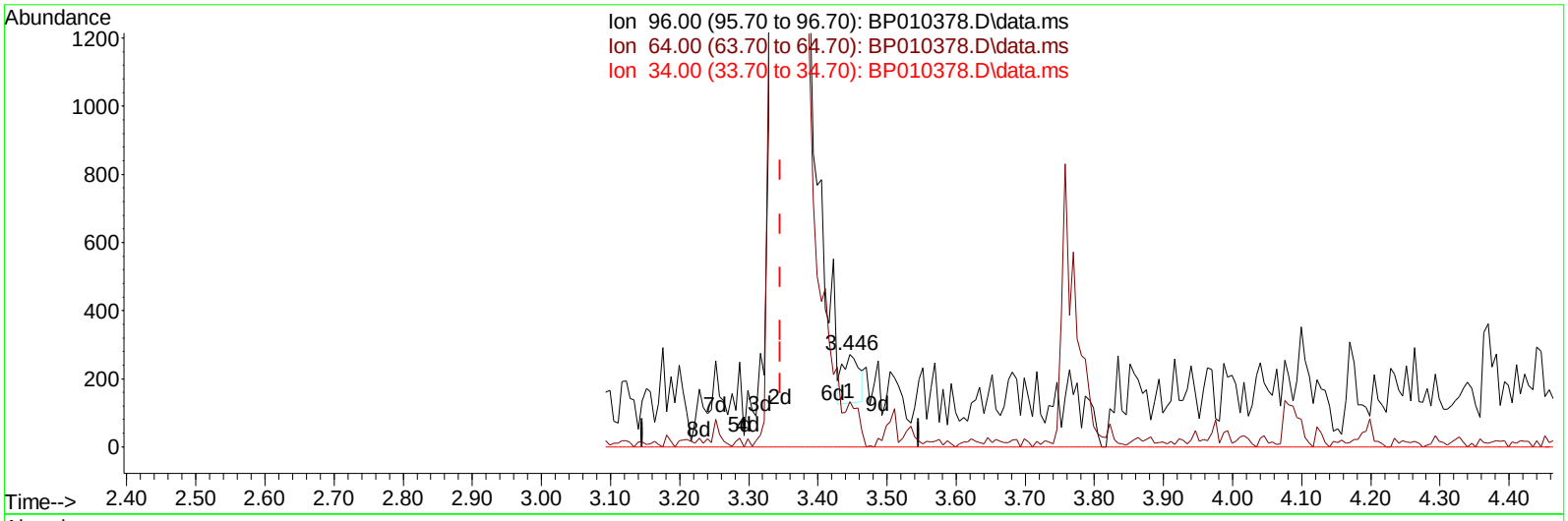
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051622\
Data File : BP010378.D
Acq On : 17 May 2022 03:41
Operator : CG/JU
Sample : PB144800BL
Misc :
ALS Vial : 26 Sample Multiplier: 1

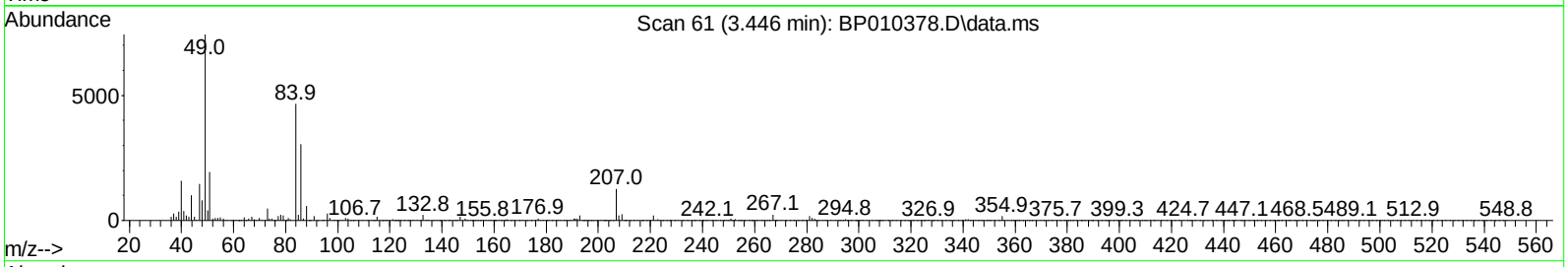
Quant Time: May 17 05:02:38 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP051322.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri May 13 15:04:31 2022
Response via : Initial Calibration

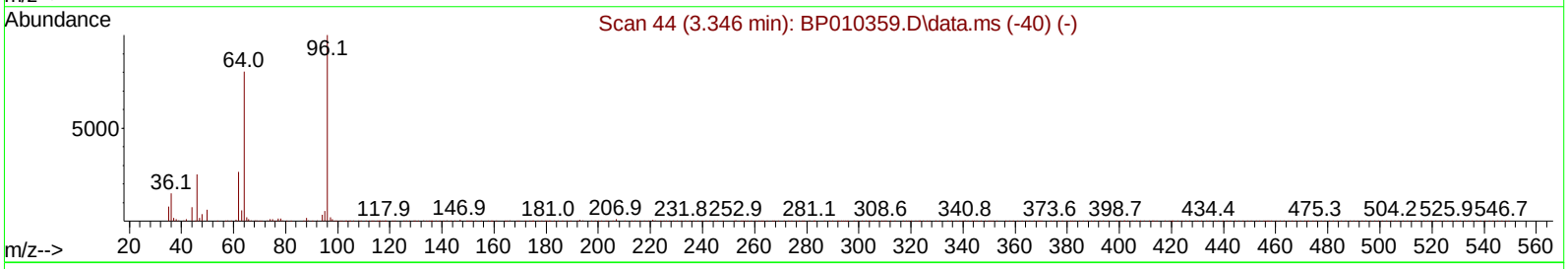
Instrument :
BNA_P
ClientSampleId :
SBLK800

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 05/17/2022
Supervised By :mohammad ahmed 05/18/2022







TIC: BP010378.D\data.ms

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

3.446min (+ 0.100) 0.06 ng/uL

response 240

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	48.53
34.00	0.00	0.00
0.00	0.00	0.00

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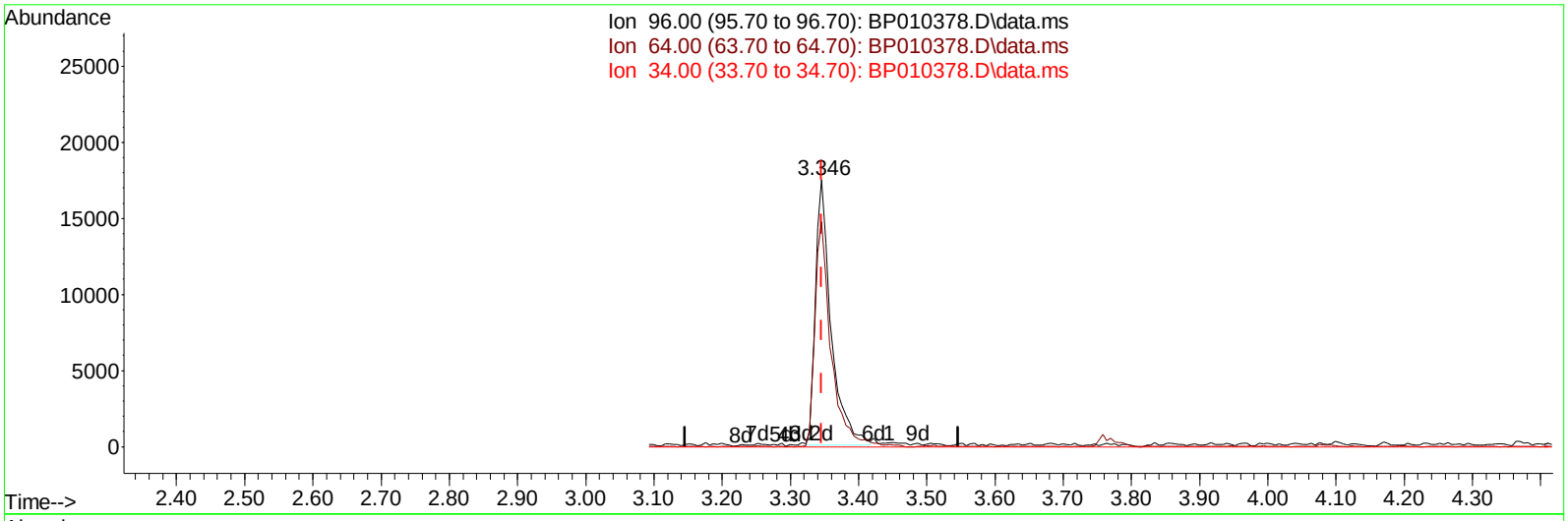

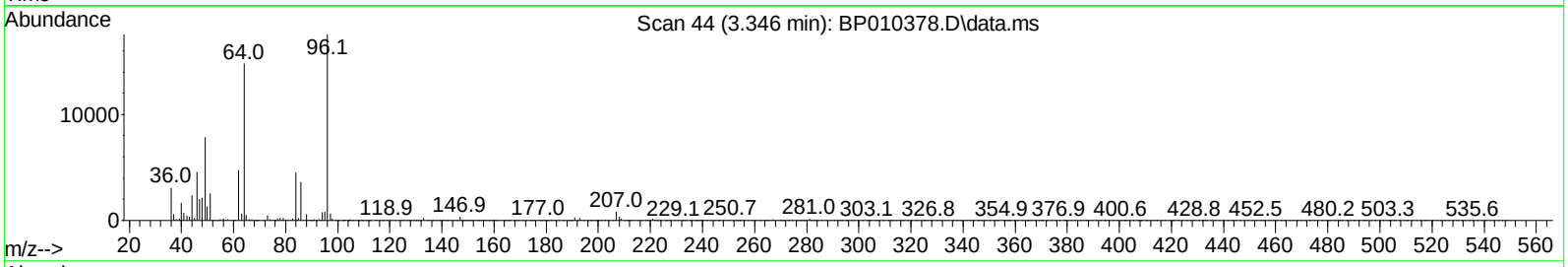

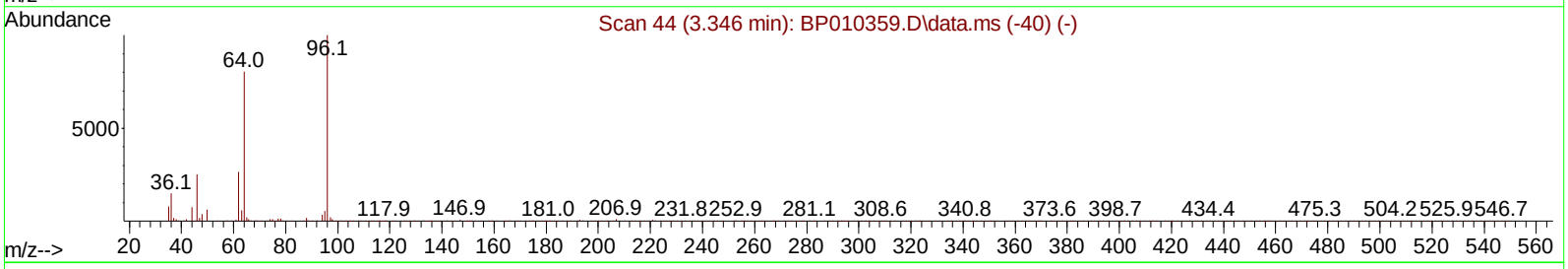

TIC: BP010378.D\data.ms

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

3.346min (0.000) 6.47 ng/uL m

response 27720

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	84.29#
34.00	0.00	0.00
0.00	0.00	0.00

Quantitation Report (Qedit)

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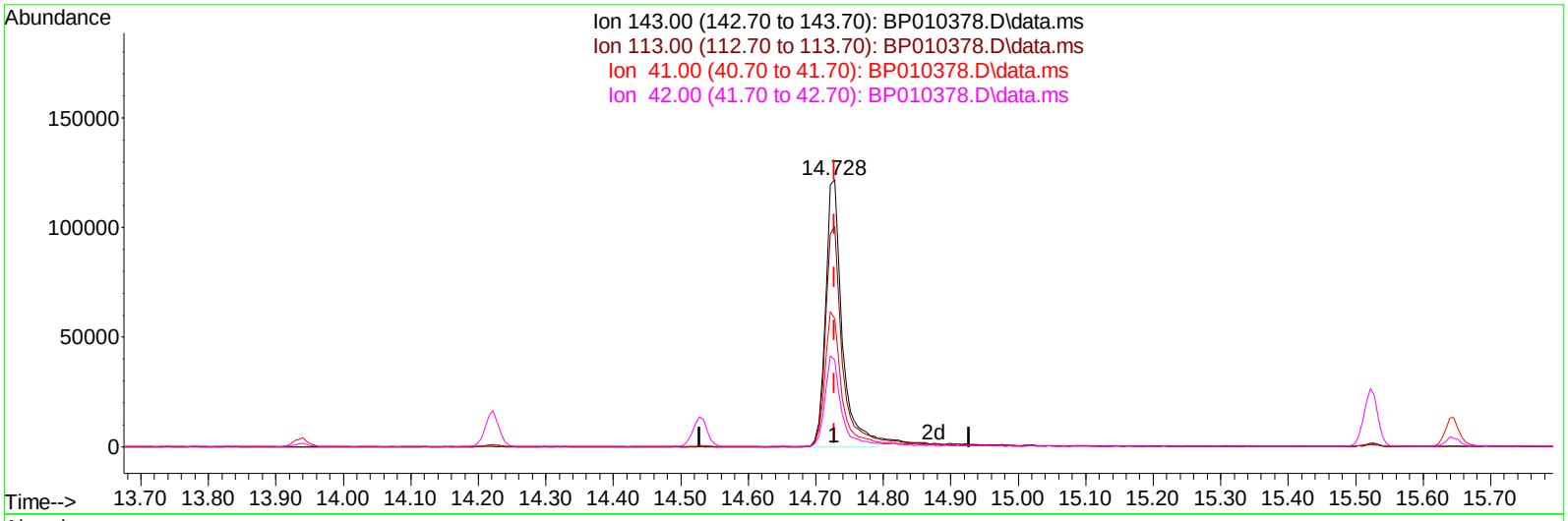

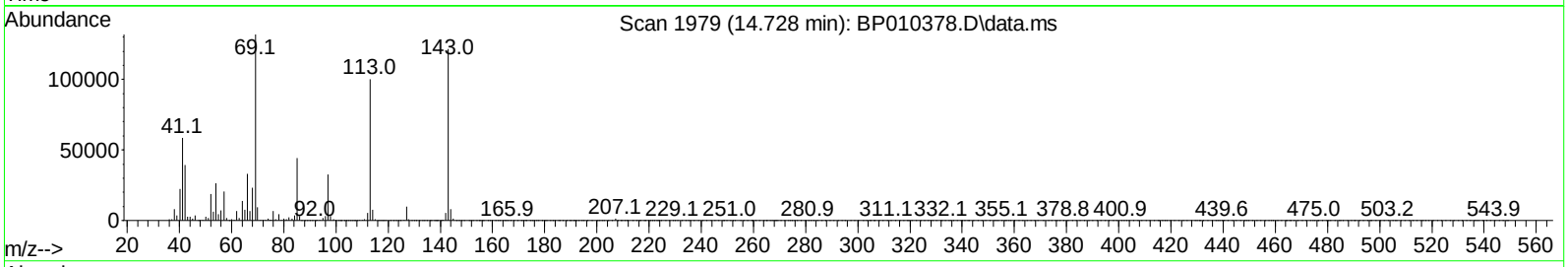

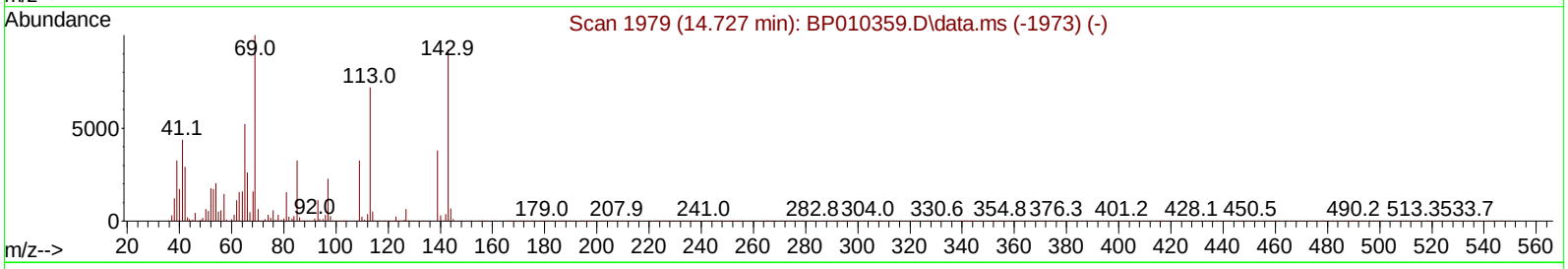

TIC: BP010378.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.728min (+ 0.000) 29.24 ng/ul

response 207793

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	132.70	82.43#
41.00	23.20	47.97#
42.00	18.00	32.36#

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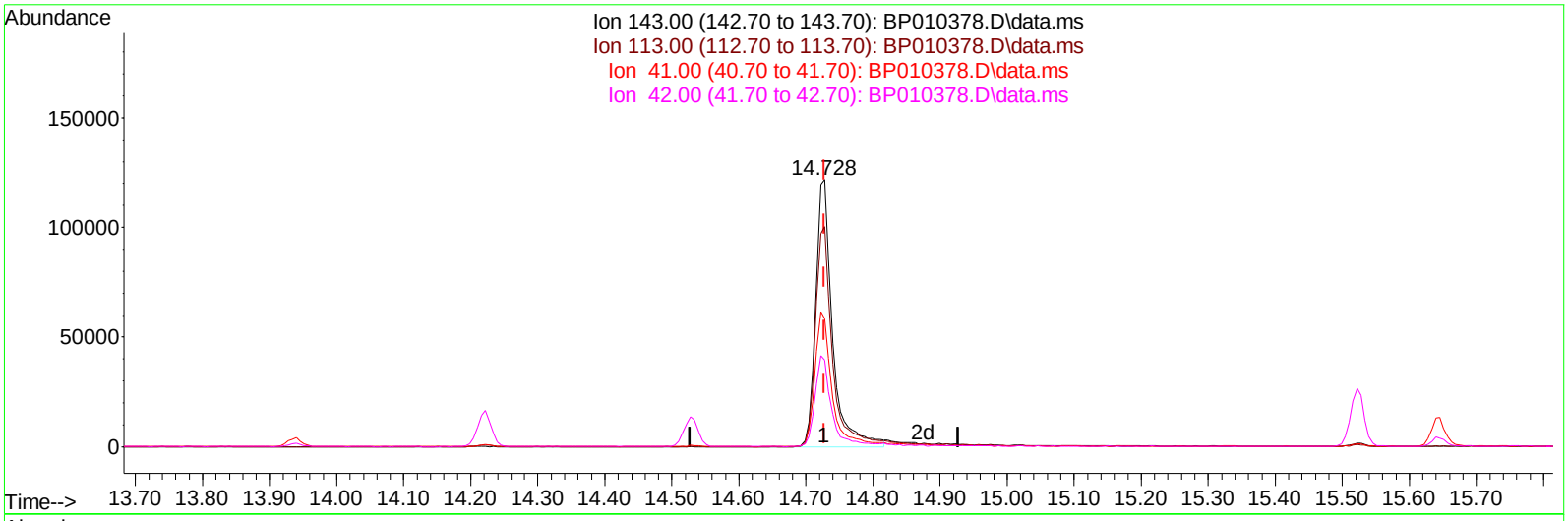

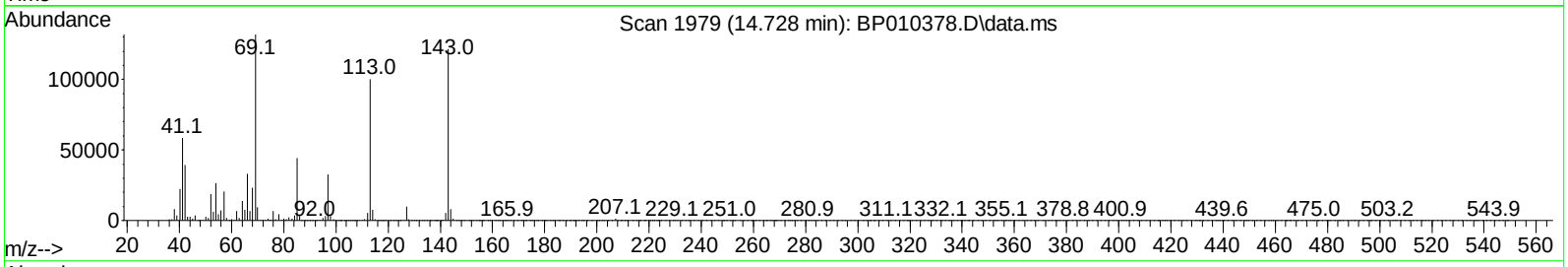

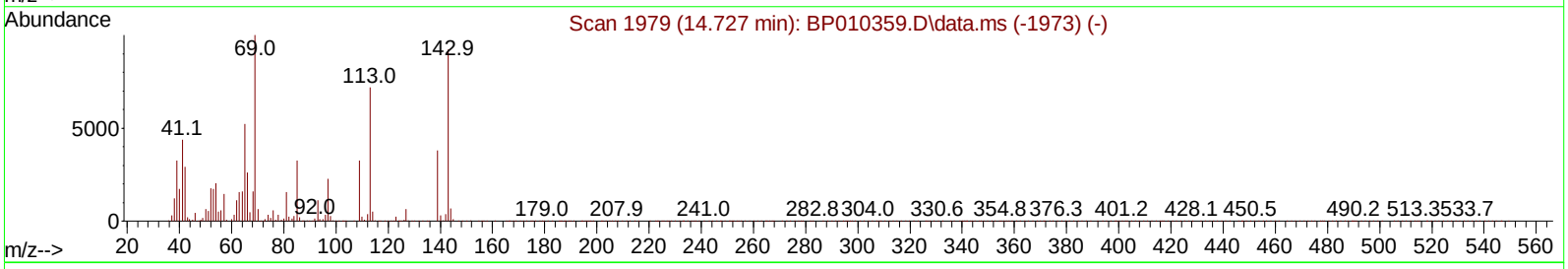

TIC: BP010378.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.728min (+ 0.000) 29.92 ng/ul m

response 212607

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	132.70	82.43#
41.00	23.20	47.97#
42.00	18.00	32.36#

Quantitation Report (QT Reviewed)

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.899	152	182574	20.000	ng/ul	0.00
20) Naphthalene-d8	10.699	136	798996	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.528	164	541941	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.281	188	1188517	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.363	240	1133993	20.000	ng/ul	# 0.00
88) Perylene-d12	23.792	264	912787	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.346	96	27720m	6.472	ng/uL	0.00
4) Pyridine-d5	3.758	84	405715	33.359	ng/ul	0.00
7) Phenol-d5	7.063	99	505935	33.016	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.228	67	352770	34.264	ng/ul	0.00
11) 2-Chlorophenol-d4	7.428	132	408288	34.820	ng/ul	0.00
15) 4-Methylphenol-d8	8.605	113	436089	33.486	ng/ul	0.00
21) Nitrobenzene-d5	9.057	128	217773	35.217	ng/ul	0.00
24) 2-Nitrophenol-d4	9.781	143	242183	36.956	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.322	165	413257	33.351	ng/ul	0.00
31) 4-Chloroaniline-d4	10.834	131	603461	32.754	ng/ul	0.00
46) Dimethylphthalate-d6	13.940	166	1472831	36.887	ng/ul	0.00
49) Acenaphthylene-d8	14.222	160	1683086	36.544	ng/ul	0.00
54) 4-Nitrophenol-d4	14.728	143	212607m	29.915	ng/ul	0.00
60) Fluorene-d10	15.522	176	1270462	36.574	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.645	200	189034	27.196	ng/ul	0.00
73) Anthracene-d10	17.381	188	1998786	37.140	ng/ul	0.00
81) Pyrene-d10	19.616	212	2373706	39.415	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.639	264	1748693	38.152	ng/ul	0.00

Target Compounds Qvalue

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

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