

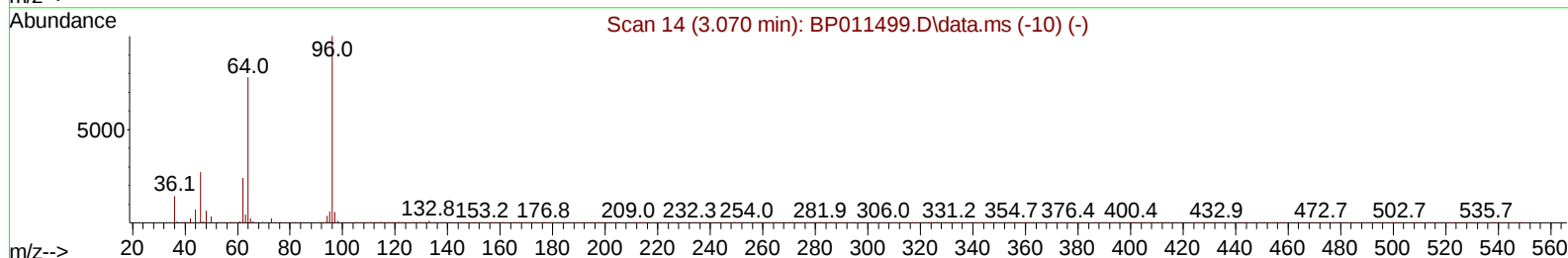
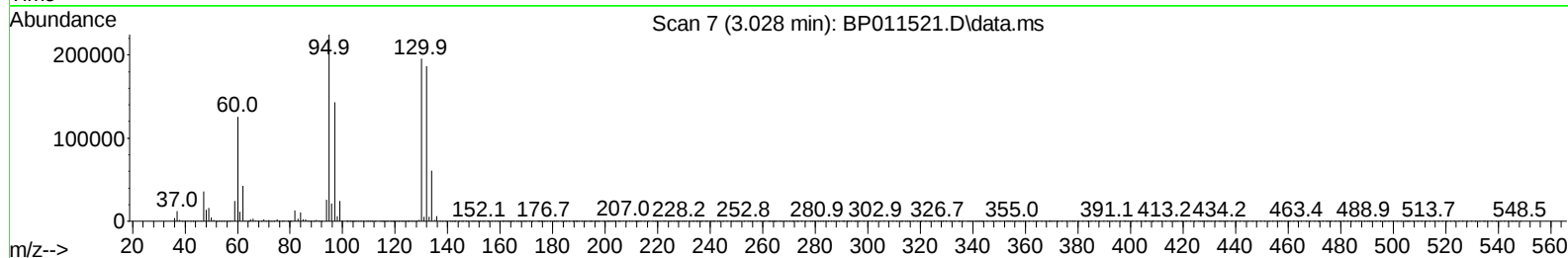
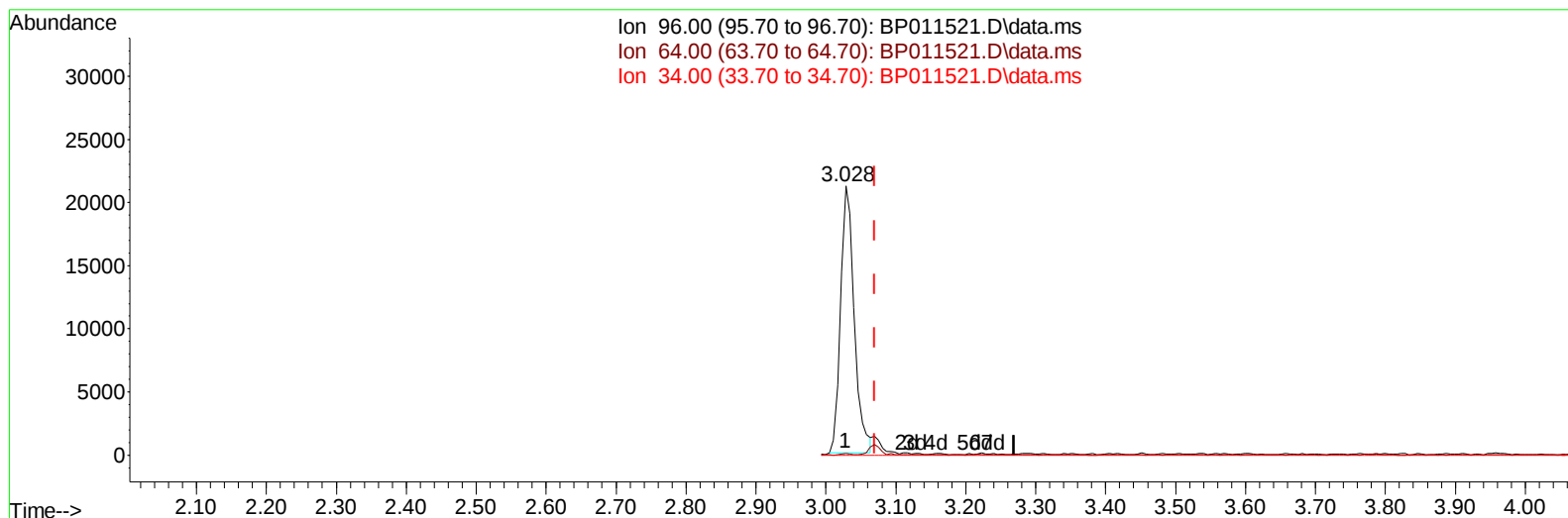
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP081922\
 Data File : BP011521.D
 Acq On : 19 Aug 2022 23:33
 Operator : CG/JU
 Sample : N4221-14DL 5X
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0KN6DL

Manual IntegrationsAPPROVED

Quant Time: Aug 20 04:42:45 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP081622.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Aug 17 01:58:37 2022
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 08/22/2022
 Supervised By :Sohil Jodhani 08/22/2022



TIC: BP011521.D\data.ms

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

3.028min (-0.042) 8.55 ng/uL

response 28929

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	74.50	0.51#
34.00	0.00	0.00
0.00	0.00	0.00

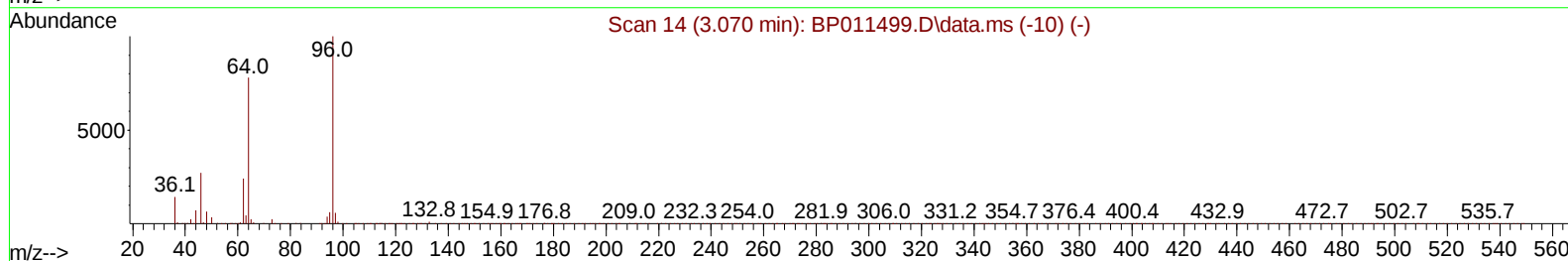
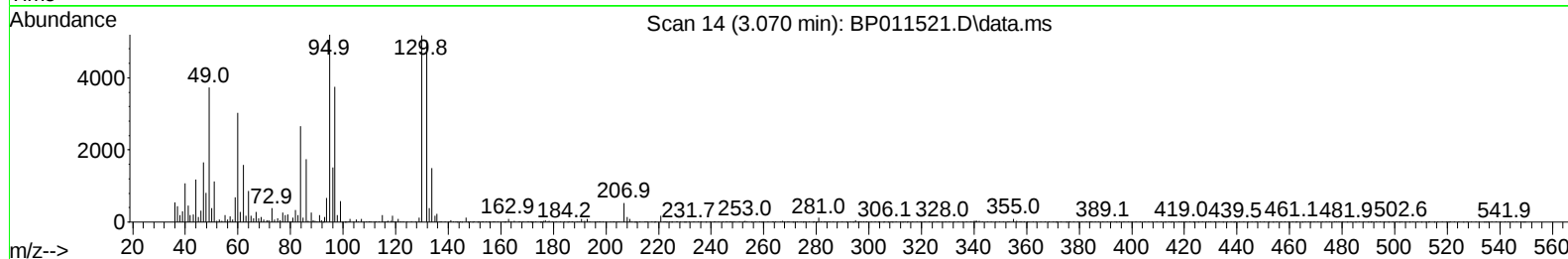
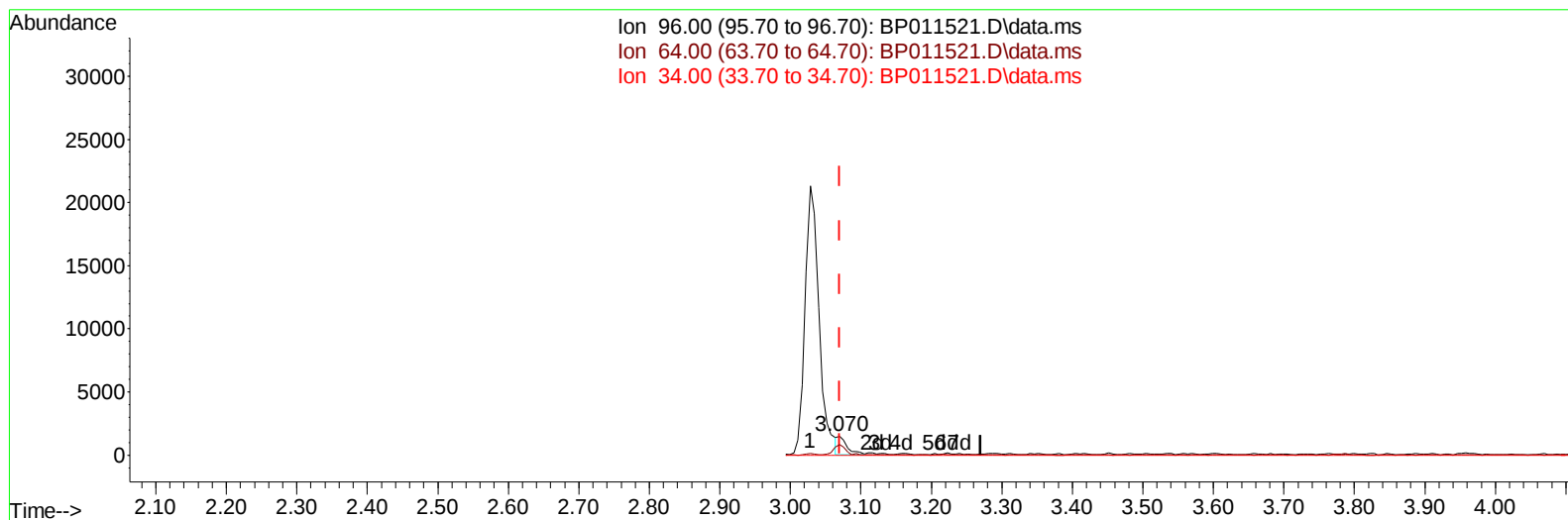
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(3) 1,4-Dioxane-d8 (S)

3.070min (-0.000) 0.41 ng/uL m

response 1371

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	74.50	56.86#
34.00	0.00	0.00
0.00	0.00	0.00

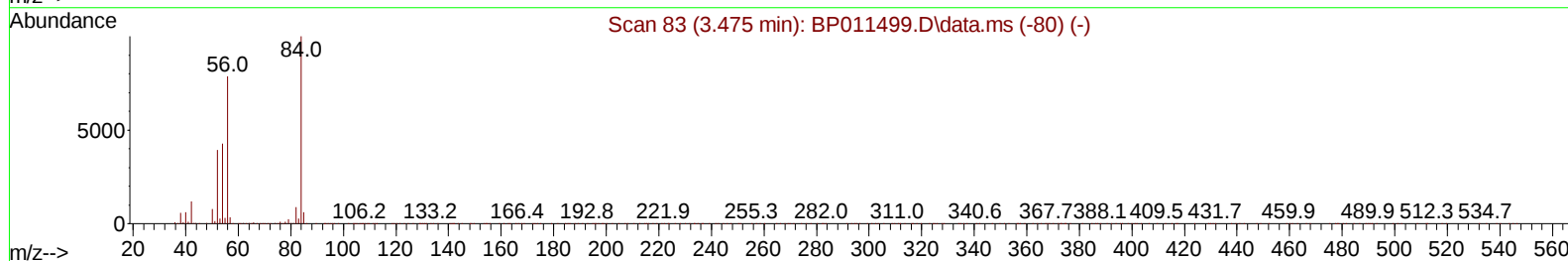
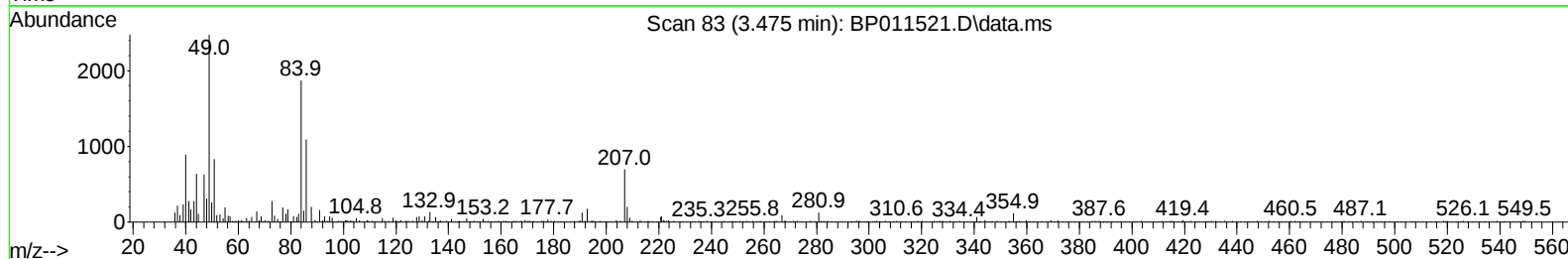
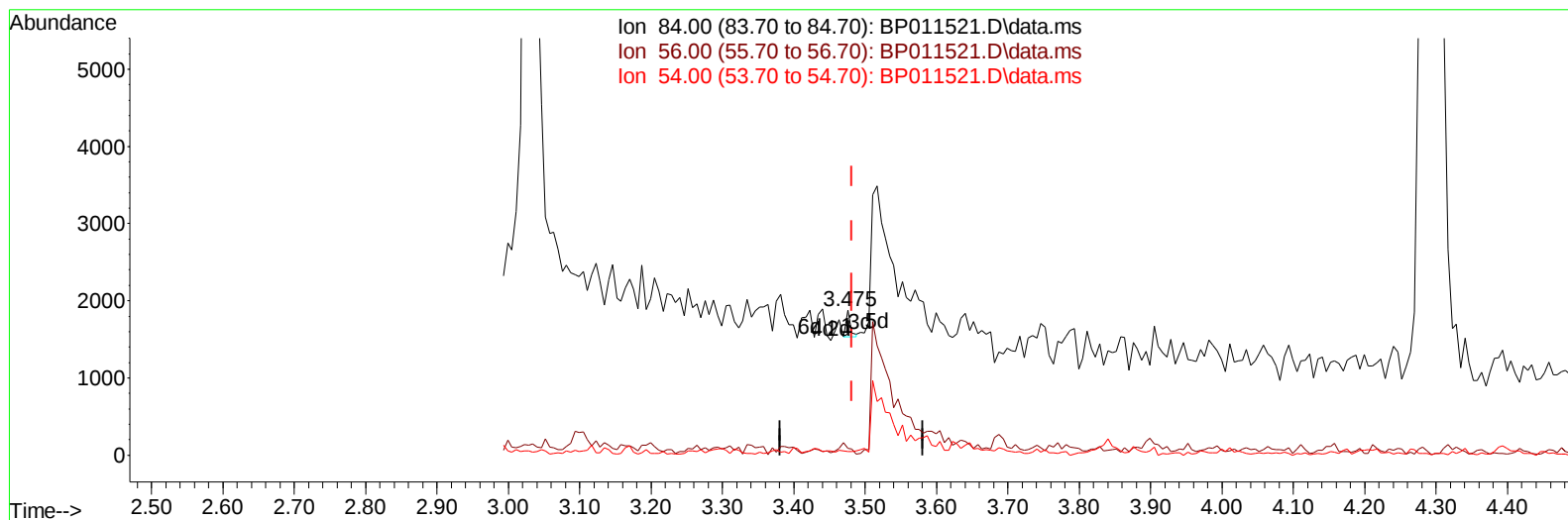
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TIC: BP011521.D\data.ms

(4) Pyridine-d5 (S)

3.475min (-0.006) 0.02 ng/u1

response 153

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	72.90	4.43#
54.00	42.10	2.56#
0.00	0.00	0.00

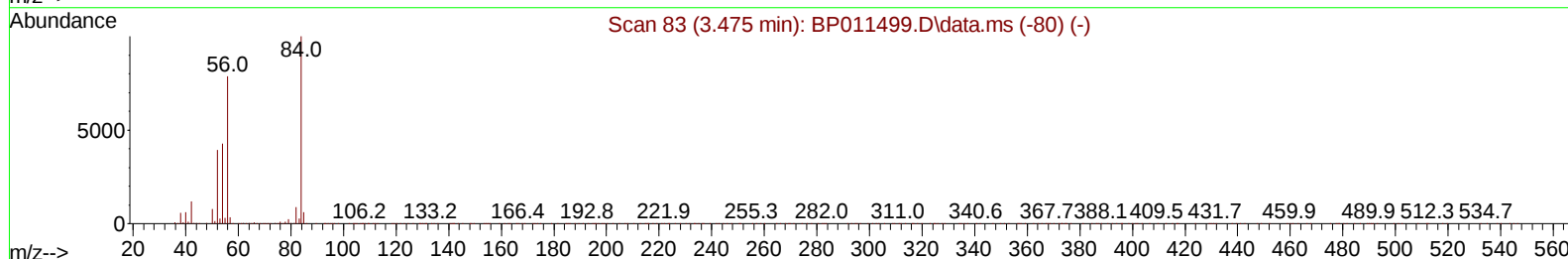
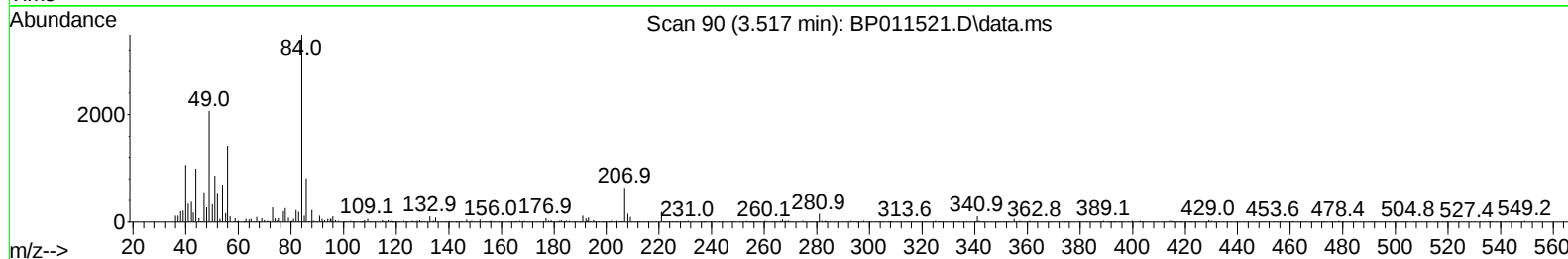
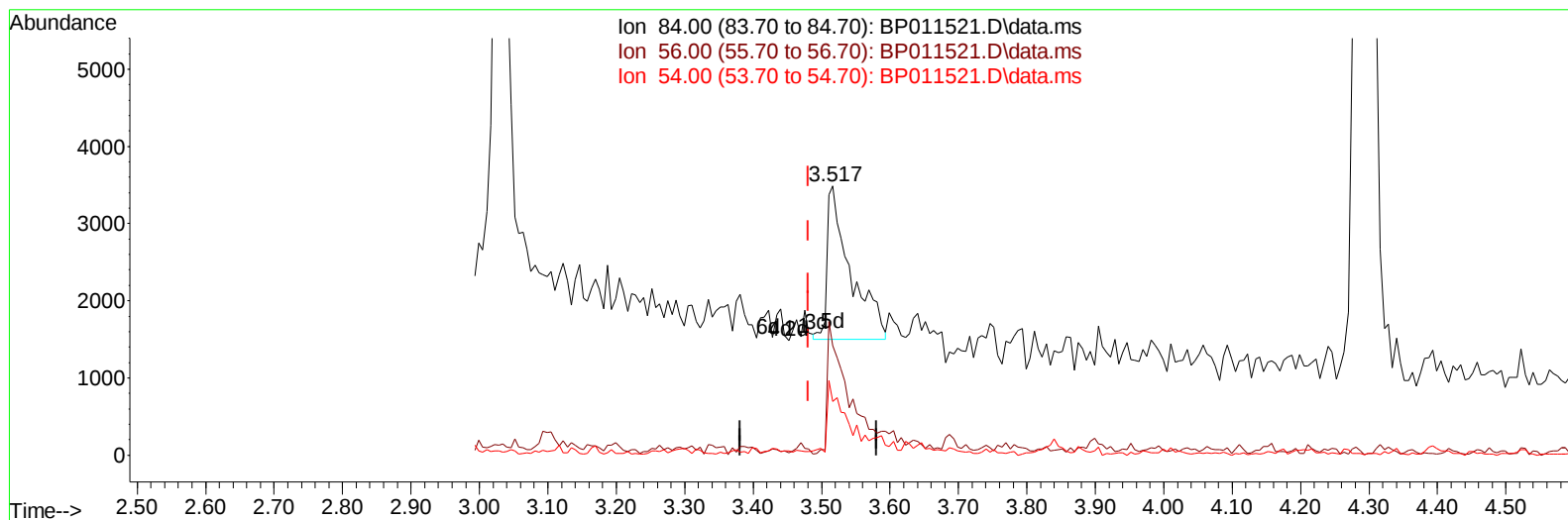
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TIC: BP011521.D\data.ms

(4) Pyridine-d5 (S)

3.517min (+ 0.036) 0.54 ng/u1 m

response 4702

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	72.90	40.59#
54.00	42.10	19.85#
0.00	0.00	0.00

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.552	152	104976	20.000	ng/ul	0.00
20) Naphthalene-d8	10.340	136	428786	20.000	ng/ul	-0.01
38) Acenaphthene-d10	14.228	164	234557	20.000	ng/ul	0.00
64) Phenanthrene-d10	16.998	188	460058	20.000	ng/ul	0.00
79) Chrysene-d12	21.122	240	454755	20.000	ng/ul	-0.01
88) Perylene-d12	23.410	264	502298	20.000	ng/ul	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.070	96	1371m	0.405	ng/uL	0.00
4) Pyridine-d5	3.517	84	4702m	0.543	ng/ul	0.04
7) Phenol-d5	6.740	99	7218	0.748	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.905	67	27022	4.312	ng/ul	0.00
11) 2-Chlorophenol-d4	7.087	132	24998	3.523	ng/ul	0.00
15) 4-Methylphenol-d8	8.275	113	13565	1.757	ng/ul	0.00
21) Nitrobenzene-d5	8.722	128	13644	4.420	ng/ul	0.00
24) 2-Nitrophenol-d4	9.434	143	12772	4.065	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	9.969	165	21089	3.637	ng/ul	0.00
31) 4-Chloroaniline-d4	10.499	131	24811	2.570	ng/ul	0.00
46) Dimethylphthalate-d6	13.651	166	83048	5.006	ng/ul	0.00
49) Acenaphthylene-d8	13.916	160	86529	4.693	ng/ul	0.00
54) 4-Nitrophenol-d4	0.000	143	0d	0.000	ng/ul	
60) Fluorene-d10	15.228	176	69395	4.972	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	0.000	200	0d	0.000	ng/ul	
73) Anthracene-d10	17.092	188	117011	5.743	ng/ul	-0.01
81) Pyrene-d10	19.357	212	129323	5.621	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.263	264	135912	5.562	ng/ul	-0.02
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.099	88	70394	19.957	ng/uL	98

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

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