

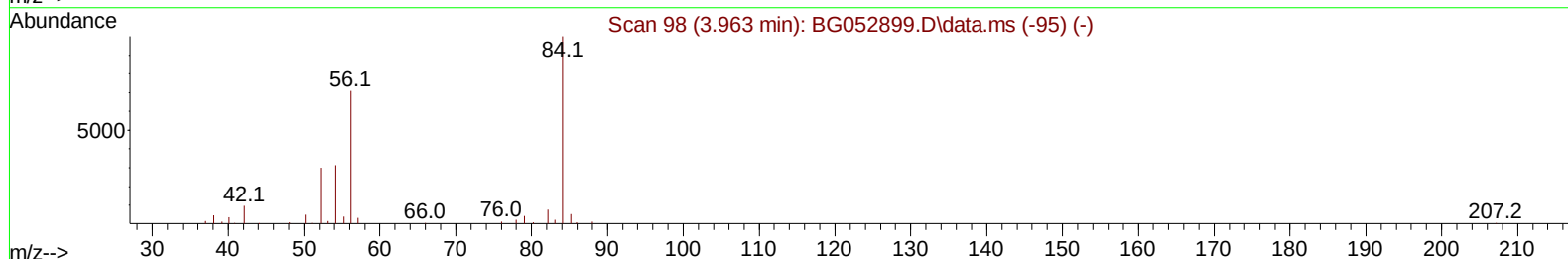
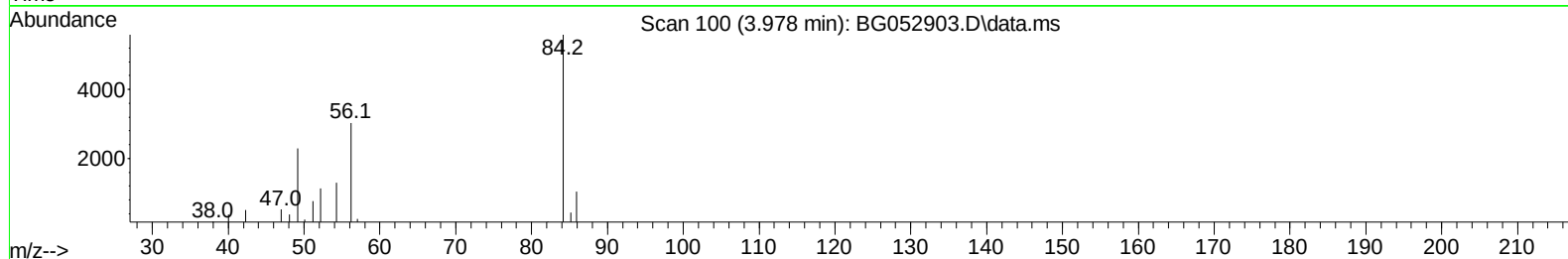
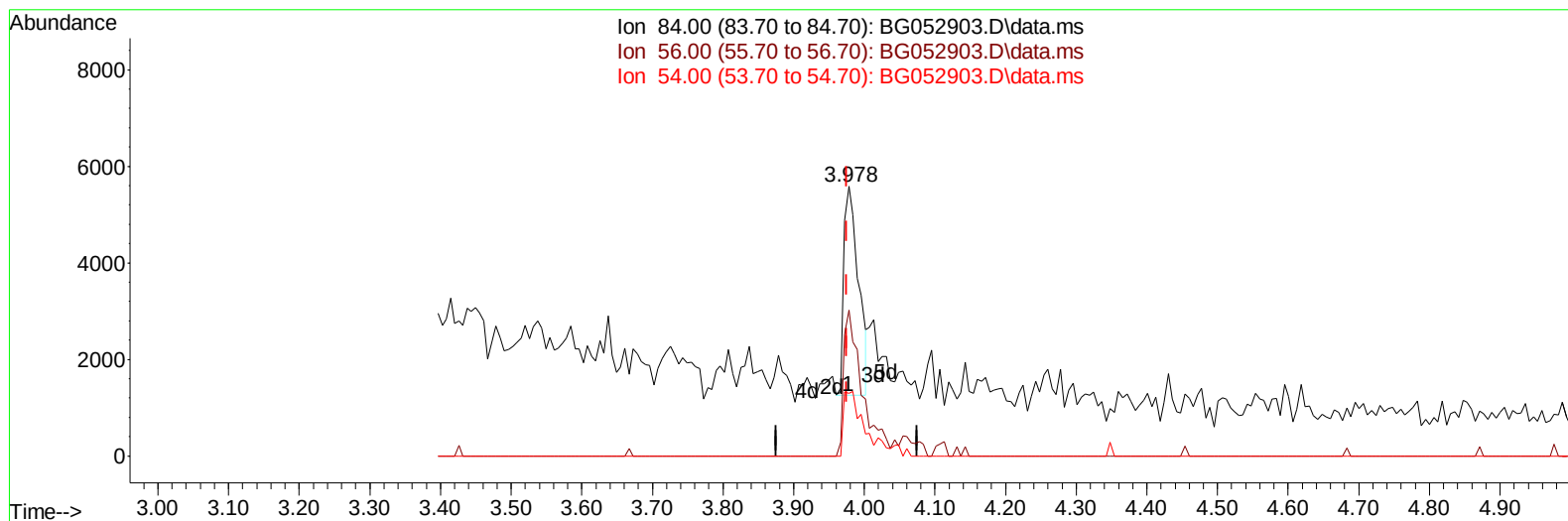
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG032822\
Data File : BG052903.D
Acq On : 28 Mar 2022 23:09
Operator : CG/JU
Sample : N2100-02
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
C0AF3

Manual IntegrationsAPPROVED

Quant Time: Mar 29 02:36:24 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG031822.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed Mar 23 13:16:32 2022
Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 03/29/2022
Supervised By :mohammad ahmed 03/29/2022



TIC: BG052903.D\data.ms

(4) Pyridine-d5 (S)

3.978min (+ 0.003) 3.71 ng/u1

response 6240

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	65.20	54.17
54.00	28.70	23.34
0.00	0.00	0.00

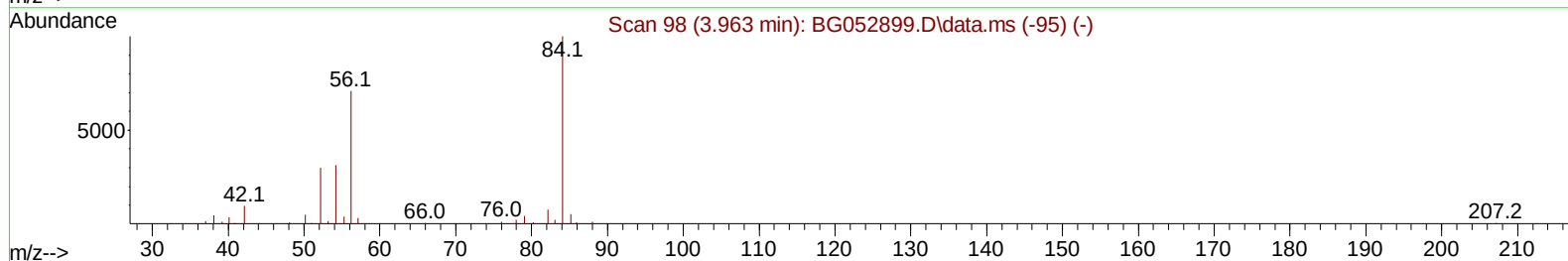
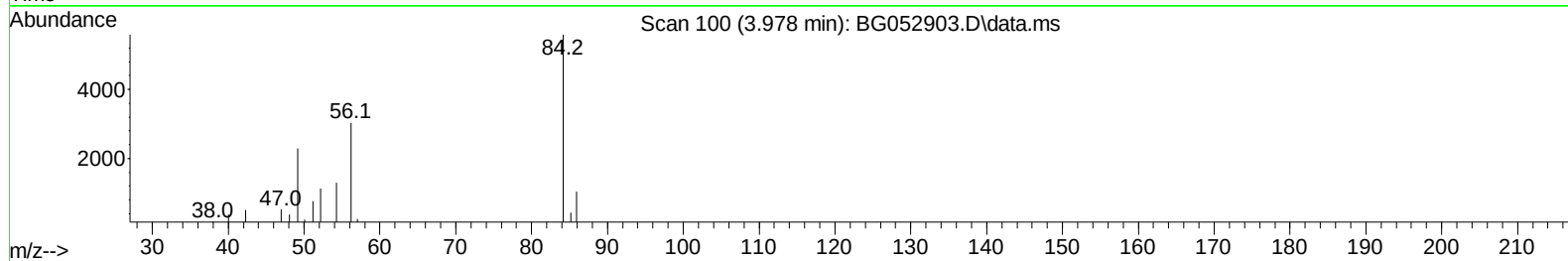
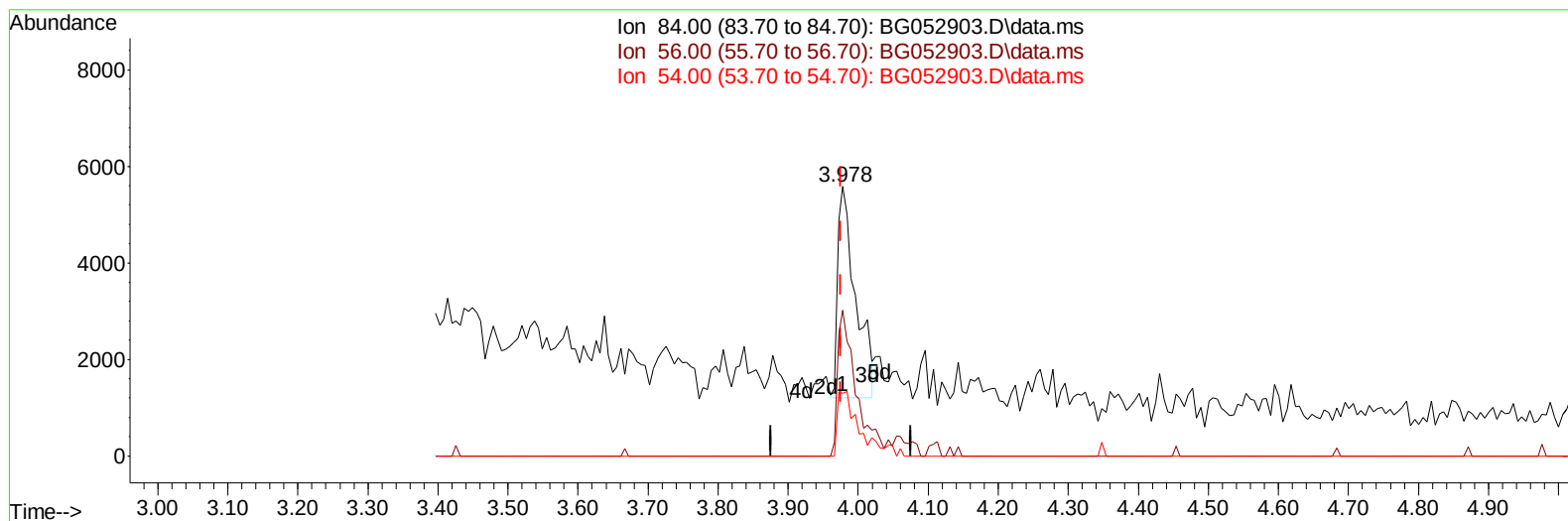
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(4) Pyridine-d5 (S)

3.978min (+ 0.003) 4.59 ng/ul m

response 7730

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	65.20	54.17
54.00	28.70	23.34
0.00	0.00	0.00

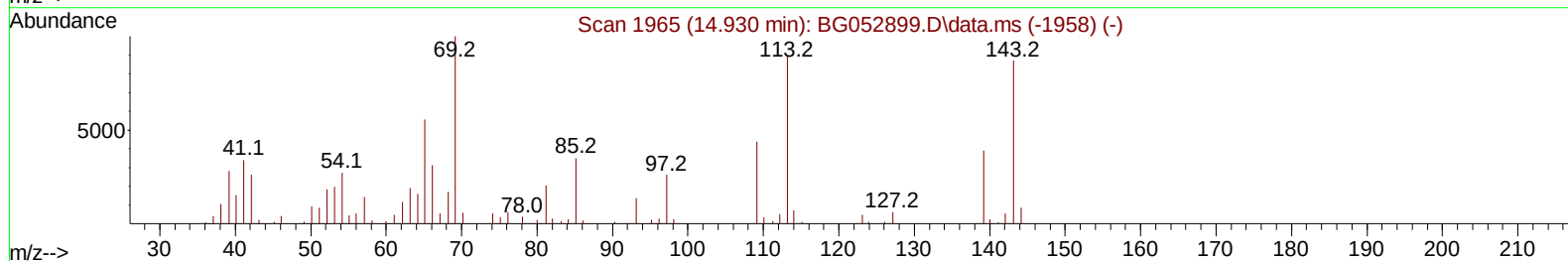
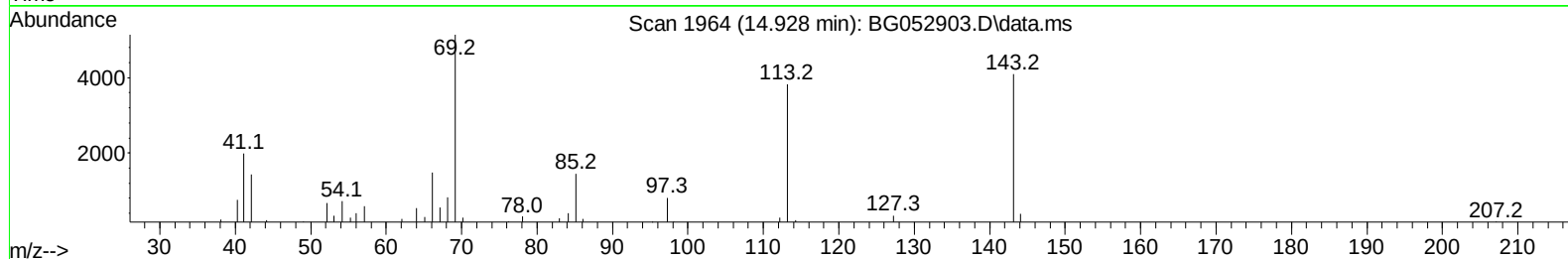
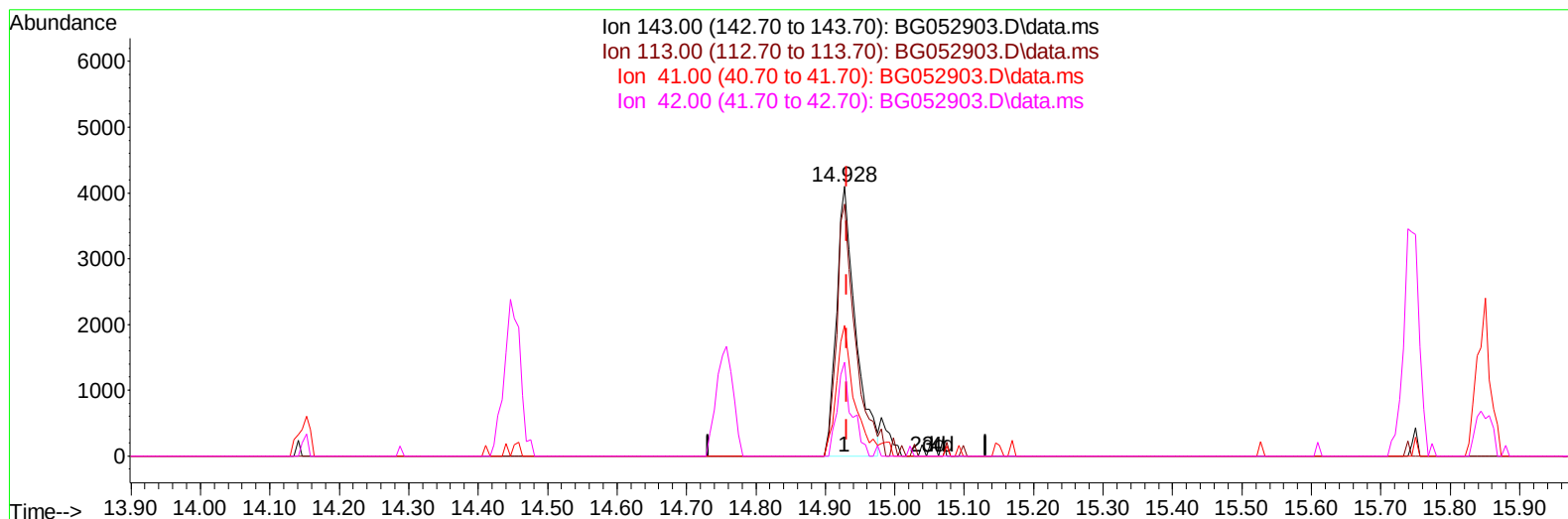
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TIC: BG052903.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.928min (-0.003) 6.81 ng/ul

response 7867

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	105.70	93.58
41.00	47.20	48.39
42.00	29.40	34.87

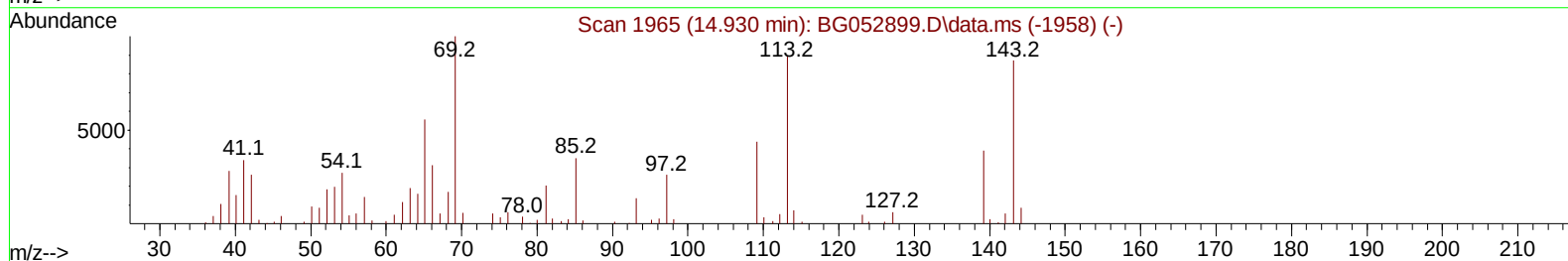
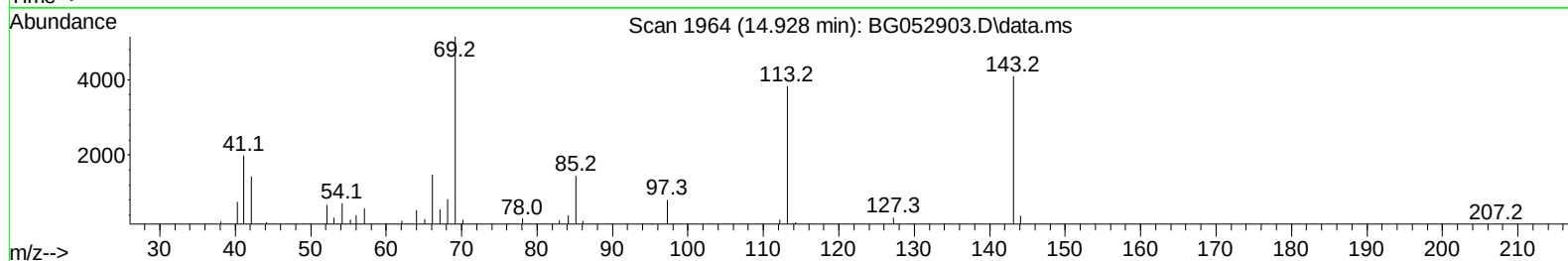
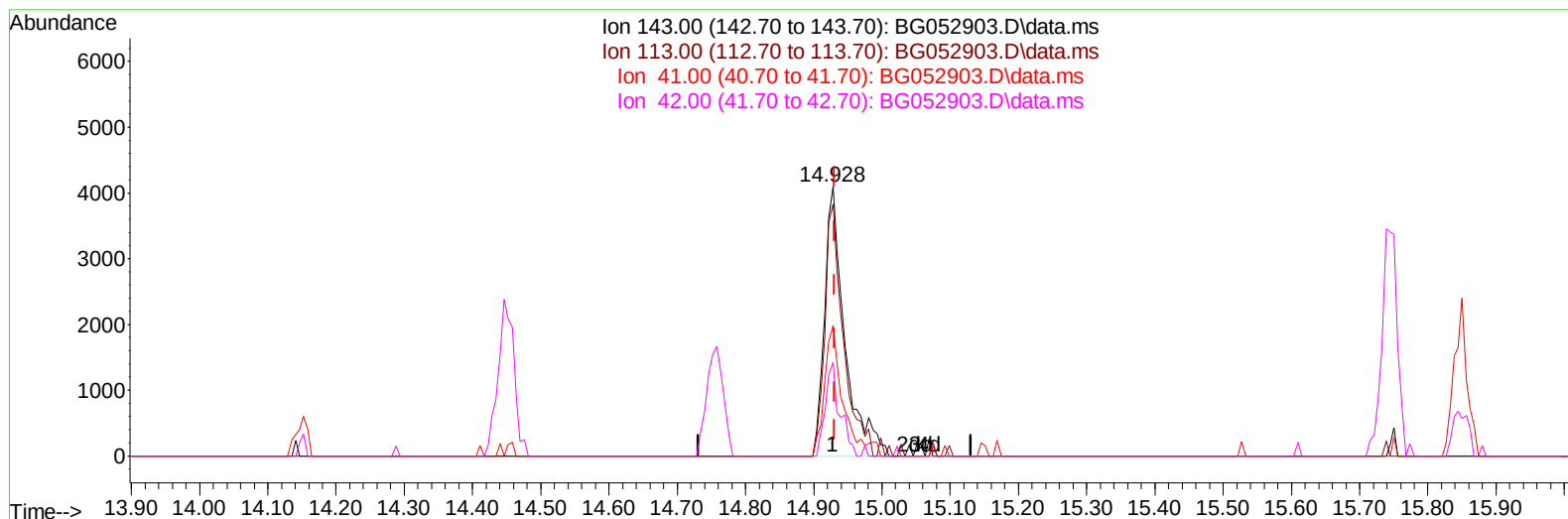
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TIC: BG052903.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.928min (-0.003) 7.32 ng/ul m

response 8450

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	105.70	93.58
41.00	47.20	48.39
42.00	29.40	34.87

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.137	152	23804	20.000	ng/ul	0.00
20) Naphthalene-d8	10.945	136	106063	20.000	ng/ul	-0.01
38) Acenaphthene-d10	14.758	164	89418	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.501	188	217779	20.000	ng/ul	0.00
79) Chrysene-d12	21.783	240	239310	20.000	ng/ul	# 0.00
88) Perylene-d12	25.102	264	238456	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.549	96	3419	5.073	ng/uL	0.00
4) Pyridine-d5	3.978	84	7730m	4.595	ng/ul	0.00
7) Phenol-d5	7.279	99	16450	7.716	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.450	67	46998	36.093	ng/ul	0.00
11) 2-Chlorophenol-d4	7.661	132	39107	27.014	ng/ul	-0.01
15) 4-Methylphenol-d8	8.830	113	28564	18.495	ng/ul	0.00
21) Nitrobenzene-d5	9.294	128	28313	37.084	ng/ul	0.00
24) 2-Nitrophenol-d4	10.017	143	30467	37.834	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.563	165	53965	30.804	ng/ul	0.00
31) 4-Chloroaniline-d4	11.074	131	61337	26.491	ng/ul	0.00
46) Dimethylphthalate-d6	14.153	166	239070	34.802	ng/ul	0.00
49) Acenaphthylene-d8	14.452	160	262672	31.467	ng/ul	0.00
54) 4-Nitrophenol-d4	14.928	143	8450m	7.316	ng/ul	0.00
60) Fluorene-d10	15.745	176	210752	34.940	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.850	200	40885	36.891	ng/ul	0.00
73) Anthracene-d10	17.595	188	385435	38.007	ng/ul	0.00
81) Pyrene-d10	19.880	212	480961	35.926	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.873	264	478413	38.829	ng/ul	0.00

Target Compounds Qvalue

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

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