

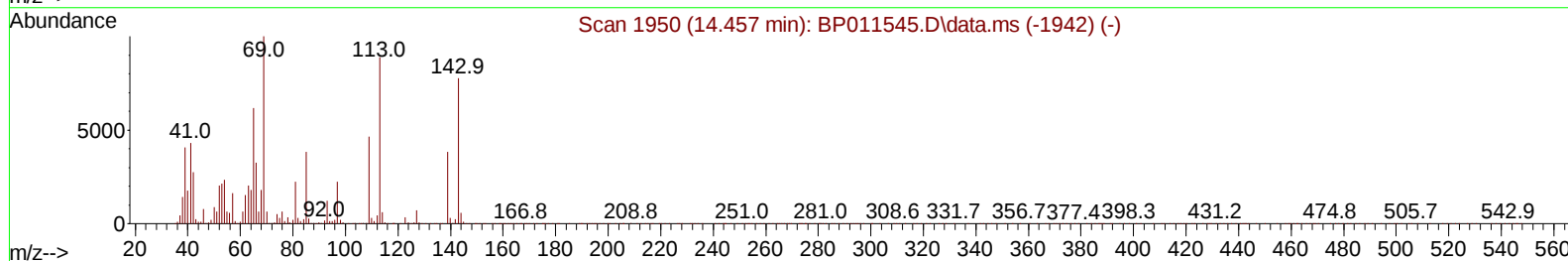
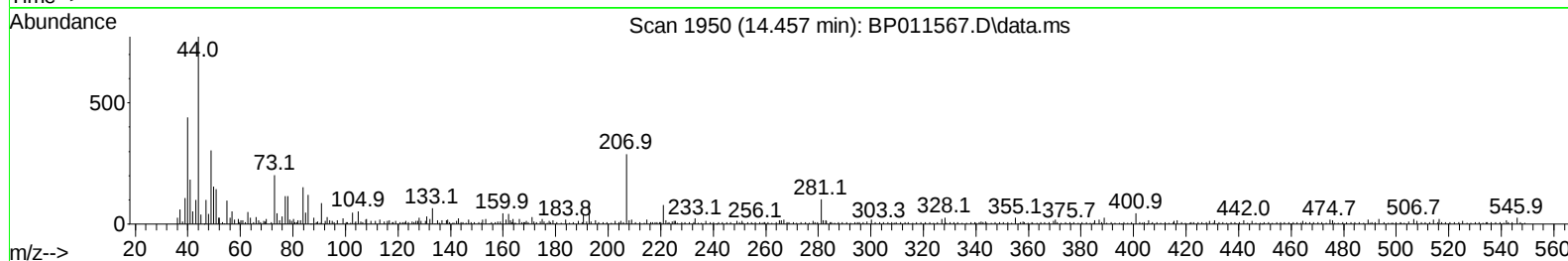
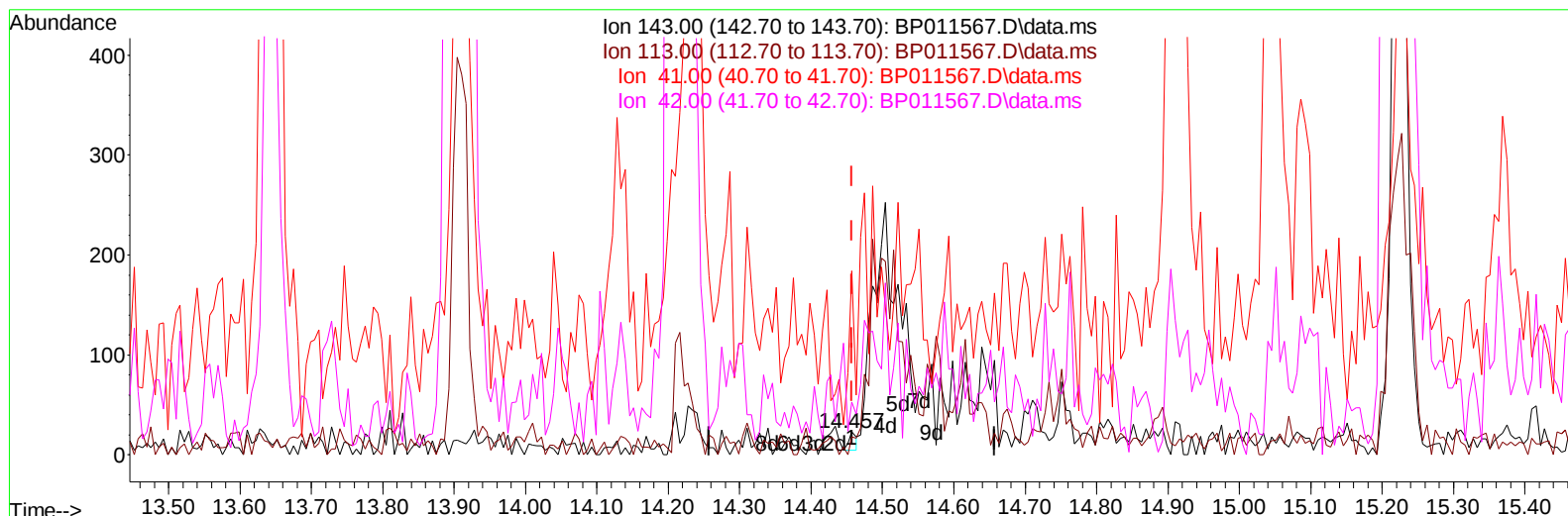
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP082422\
 Data File : BP011567.D
 Acq On : 25 Aug 2022 17:52
 Operator : CG/JU
 Sample : N4349-02
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BGN93

Manual IntegrationsAPPROVED

Quant Time: Aug 26 01:06:55 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP081622.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Aug 25 02:14:54 2022
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 08/26/2022
 Supervised By :mohammad ahmed 08/27/2022



TIC: BP011567.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.457min (+ 0.000) 0.01 ng/u1

response 13

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	109.60	78.26#
41.00	52.50	800.00#
42.00	35.70	230.43#

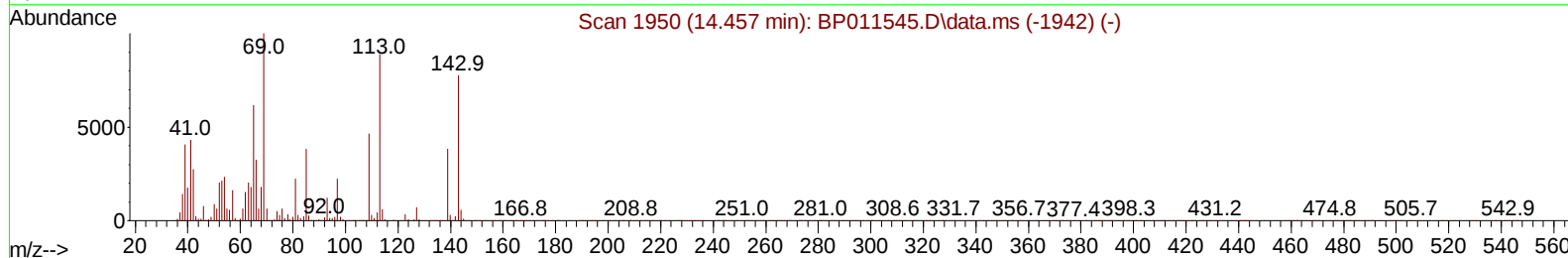
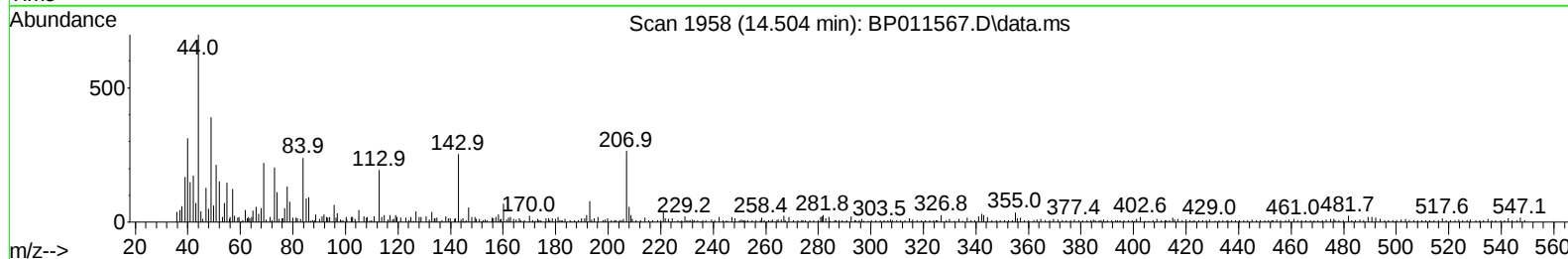
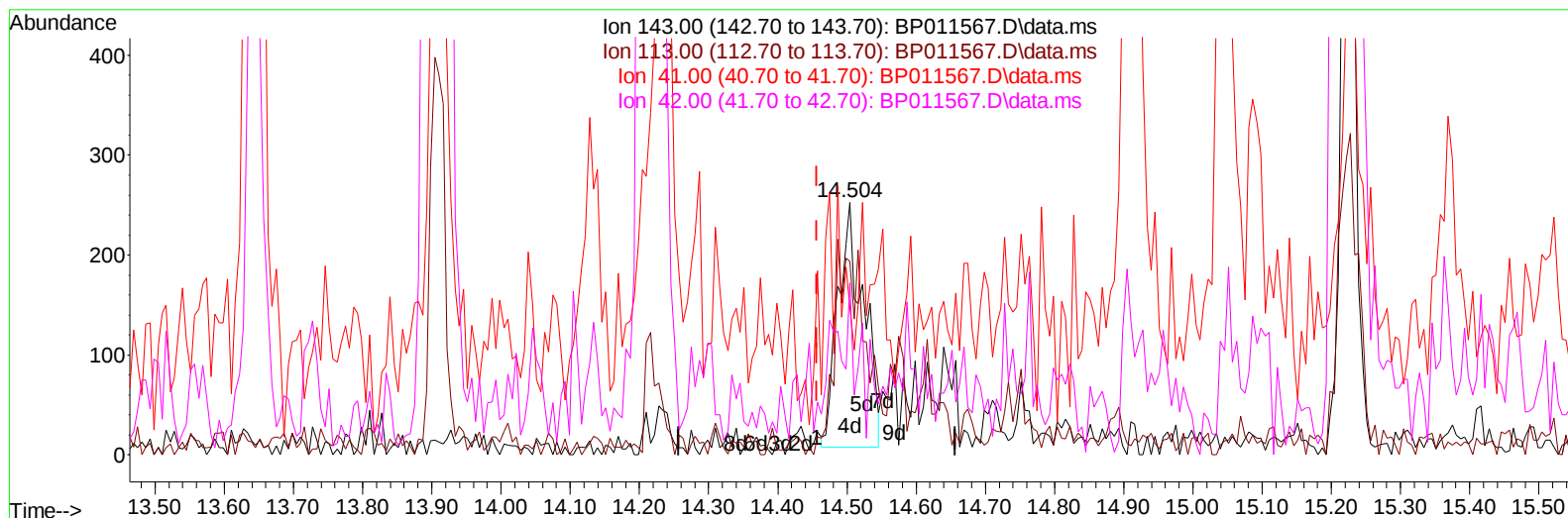
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(54) 4-Nitrophenol-d4 (S)

14.504min (+ 0.047) 0.28 ng/ul m

response 626

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	109.60	76.68#
41.00	52.50	58.50
42.00	35.70	67.98#

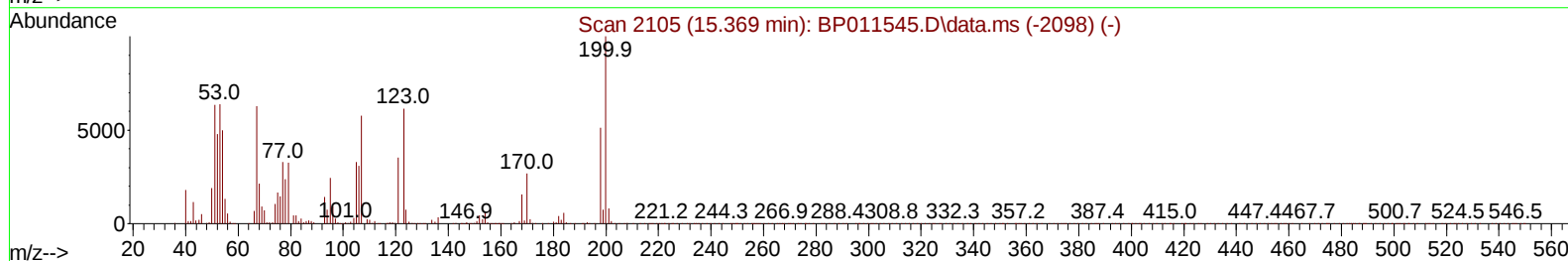
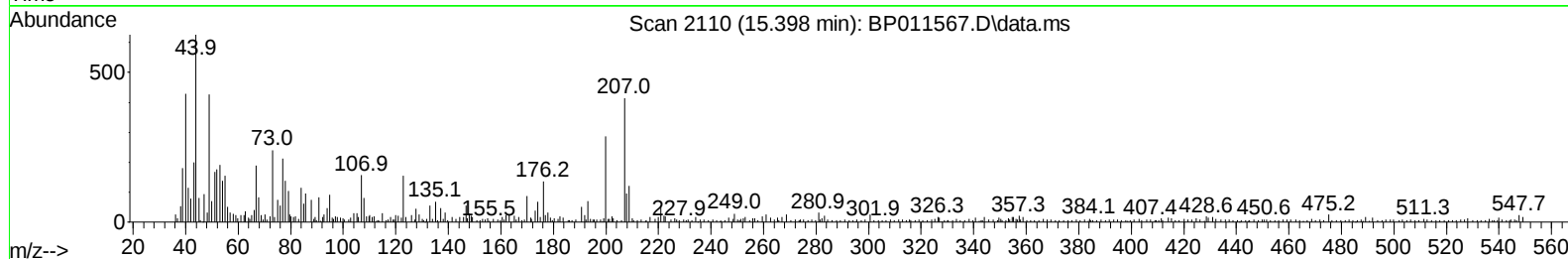
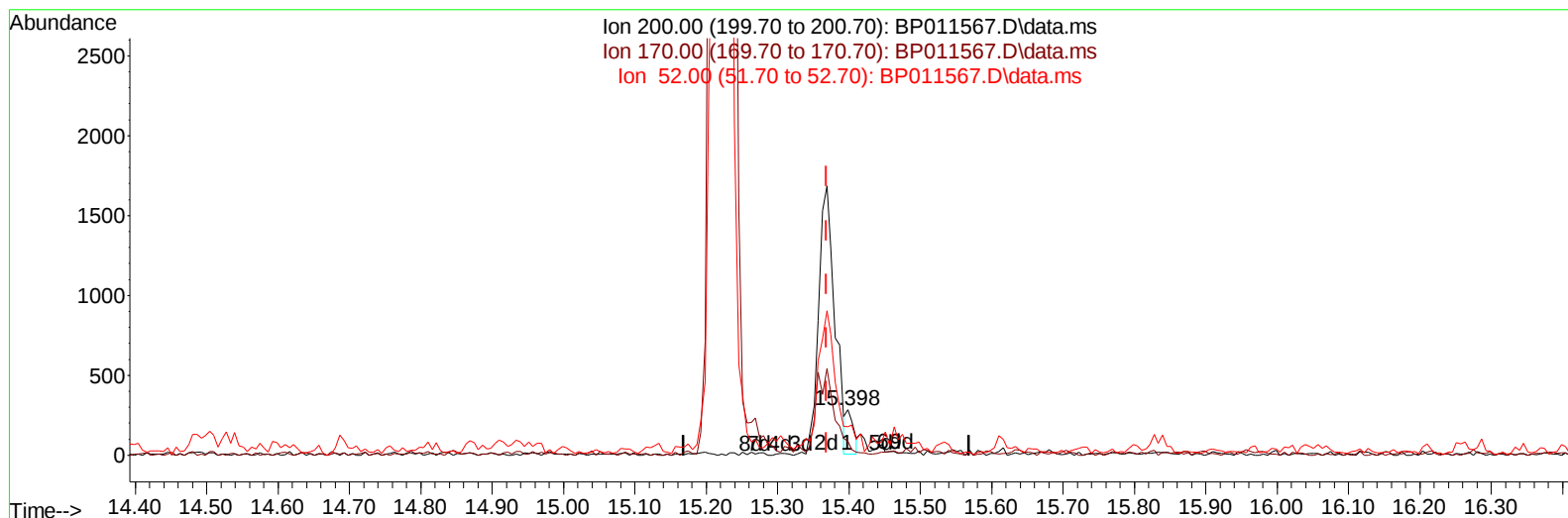
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TIC: BP011567.D\data.ms

(65) 4,6-Dinitro-2-methylphenol-d2 (S)

15.398min (+ 0.030) 0.12 ng/ul

response 209

Ion	Exp%	Act%
200.00	100.00	100.00
170.00	26.30	30.42
52.00	67.40	61.54
0.00	0.00	0.00

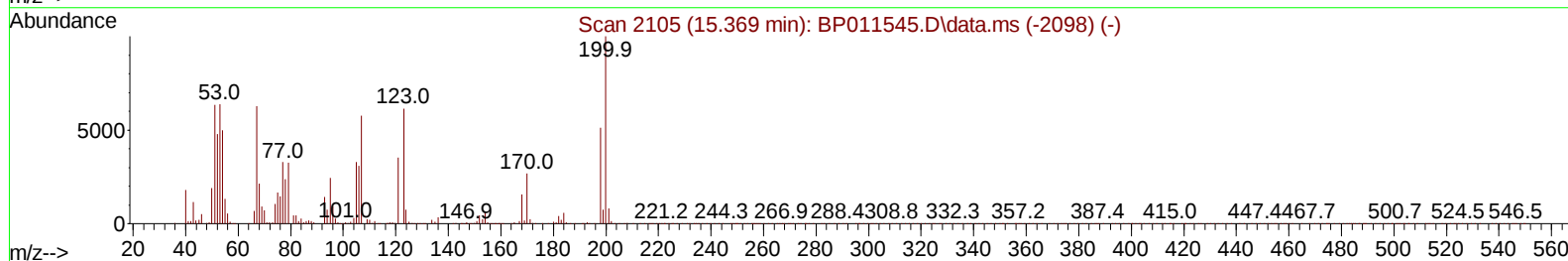
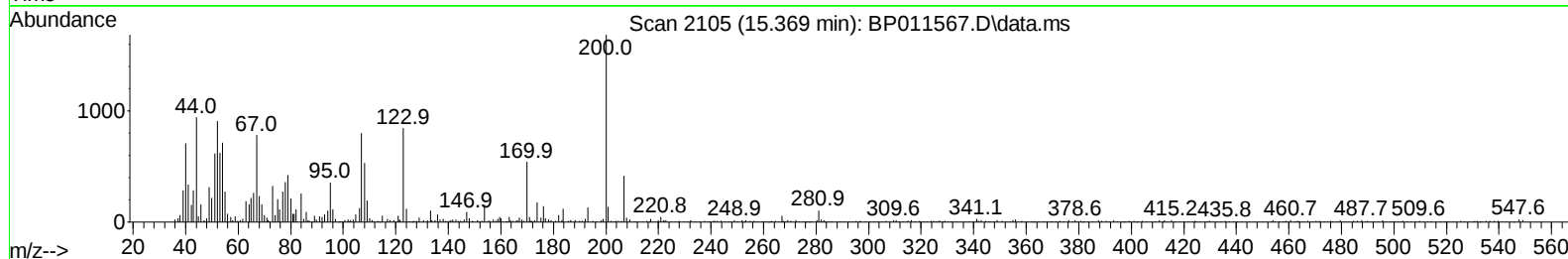
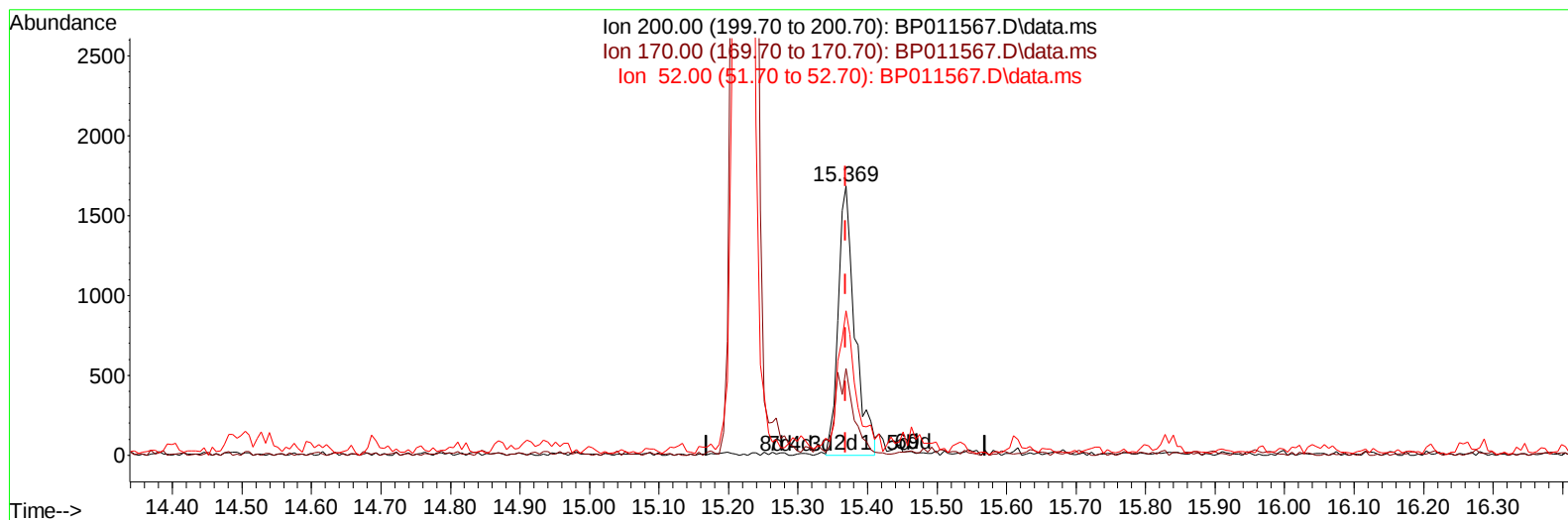
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TIC: BP011567.D\data.ms

(65) 4,6-Dinitro-2-methylphenol-d2 (S)

15.369min (+ 0.000) 1.61 ng/ul m

response 2850

Ion	Exp%	Act%
200.00	100.00	100.00
170.00	26.30	32.13#
52.00	67.40	53.74#
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	7.546	152	60592	20.000	ng/ul	0.00
20) Naphthalene-d8	10.334	136	265836	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.216	164	161460	20.000	ng/ul	0.00
64) Phenanthrene-d10	16.992	188	326739	20.000	ng/ul	0.00
79) Chrysene-d12	21.122	240	297711	20.000	ng/ul	0.00
88) Perylene-d12	23.404	264	312535	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.064	96	10271	5.260	ng/uL	0.00
4) Pyridine-d5	3.481	84	46731	9.348	ng/ul	0.00
7) Phenol-d5	6.734	99	37323	6.703	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.899	67	123827	34.230	ng/ul	0.00
11) 2-Chlorophenol-d4	7.081	132	110907	27.081	ng/ul	0.00
15) 4-Methylphenol-d8	8.269	113	71712	16.091	ng/ul	0.00
21) Nitrobenzene-d5	8.710	128	72124	37.686	ng/ul	0.00
24) 2-Nitrophenol-d4	9.428	143	38051	19.536	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	9.963	165	110271	30.671	ng/ul	0.00
31) 4-Chloroaniline-d4	10.487	131	200166	33.445	ng/ul	0.00
46) Dimethylphthalate-d6	13.646	166	435881	38.169	ng/ul	0.00
49) Acenaphthylene-d8	13.910	160	475853	37.492	ng/ul	0.00
54) 4-Nitrophenol-d4	14.504	143	626m	0.285	ng/ul	0.05
60) Fluorene-d10	15.222	176	359989	37.470	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.369	200	2850m	1.614	ng/ul	0.00
73) Anthracene-d10	17.092	188	621689	42.961	ng/ul	0.00
81) Pyrene-d10	19.357	212	666871	44.273	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.257	264	636121	41.842	ng/ul	0.00

Target Compounds Qvalue

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

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