

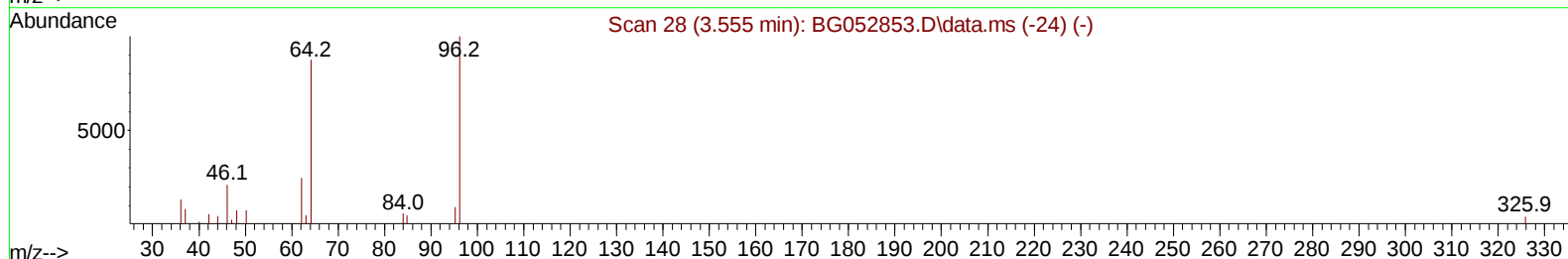
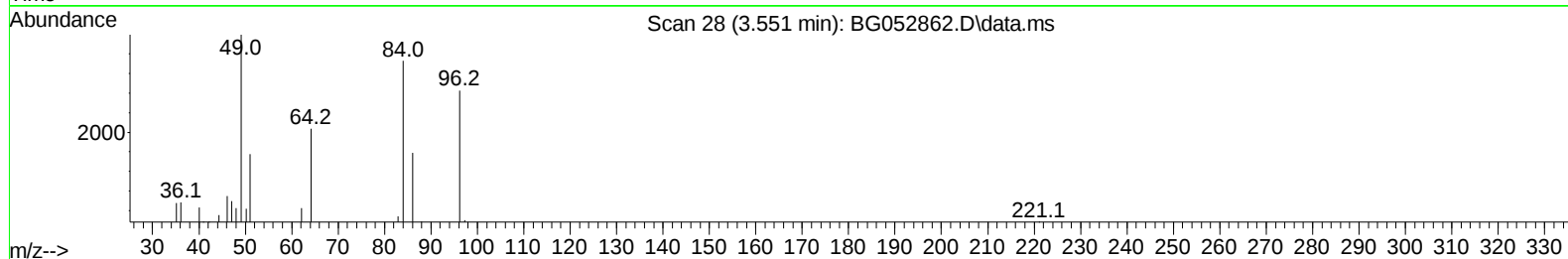
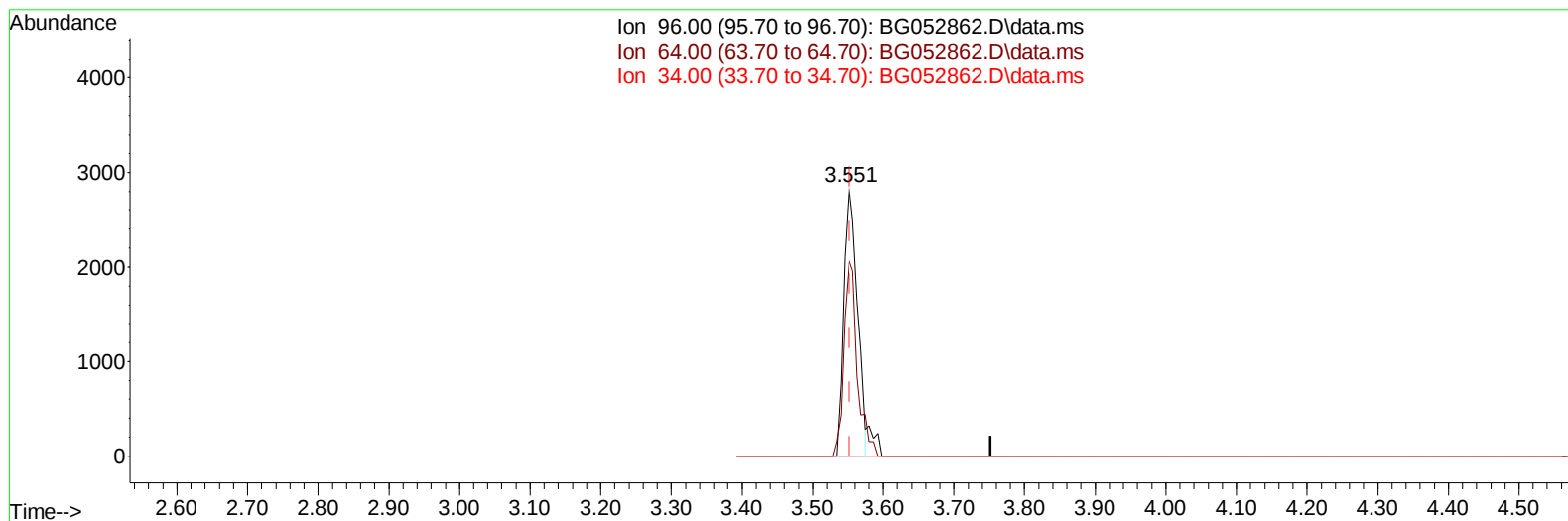
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG032522\
 Data File : BG052862.D
 Acq On : 26 Mar 2022 13:28
 Operator : CG/JU
 Sample : N2068-05
 Misc :
 ALS Vial : 40 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 YAZ10

Manual IntegrationsAPPROVED

Quant Time: Mar 28 00:59:00 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 03/28/2022
 Supervised By :mohammad ahmed 03/29/2022



TIC: BG052862.D\data.ms

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

3.551min (-0.001) 4.87 ng/uL

response 3966

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	83.40	72.78
34.00	0.00	0.00
0.00	0.00	0.00

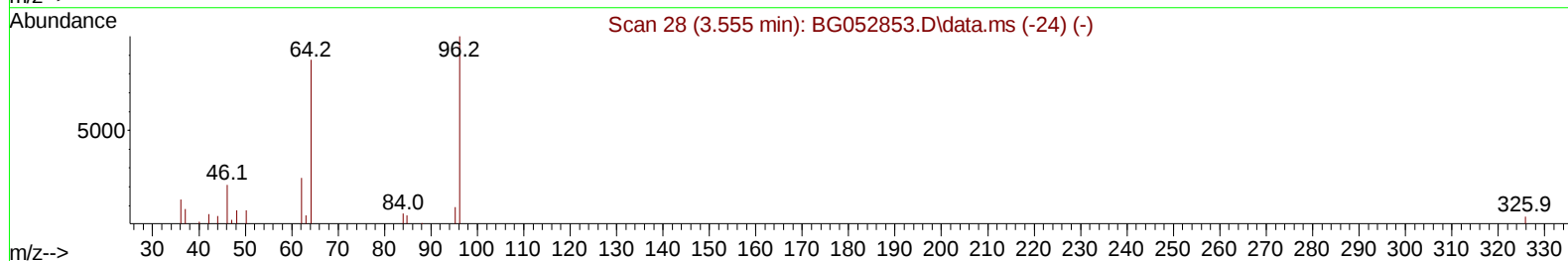
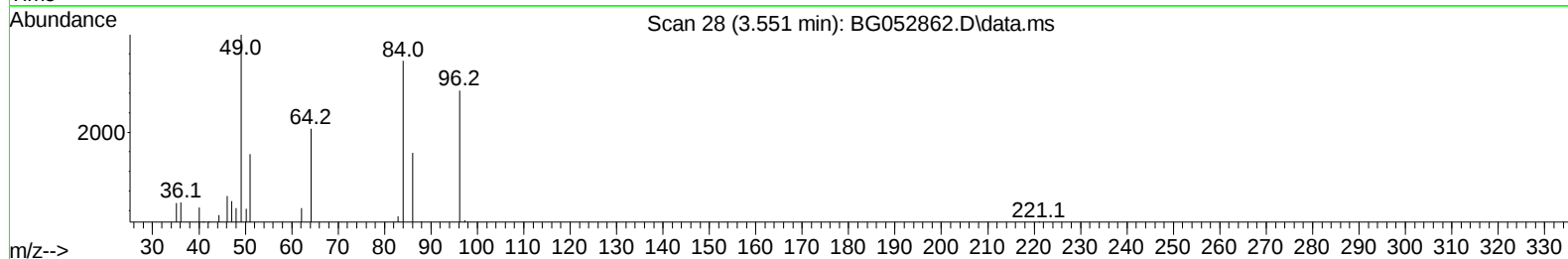
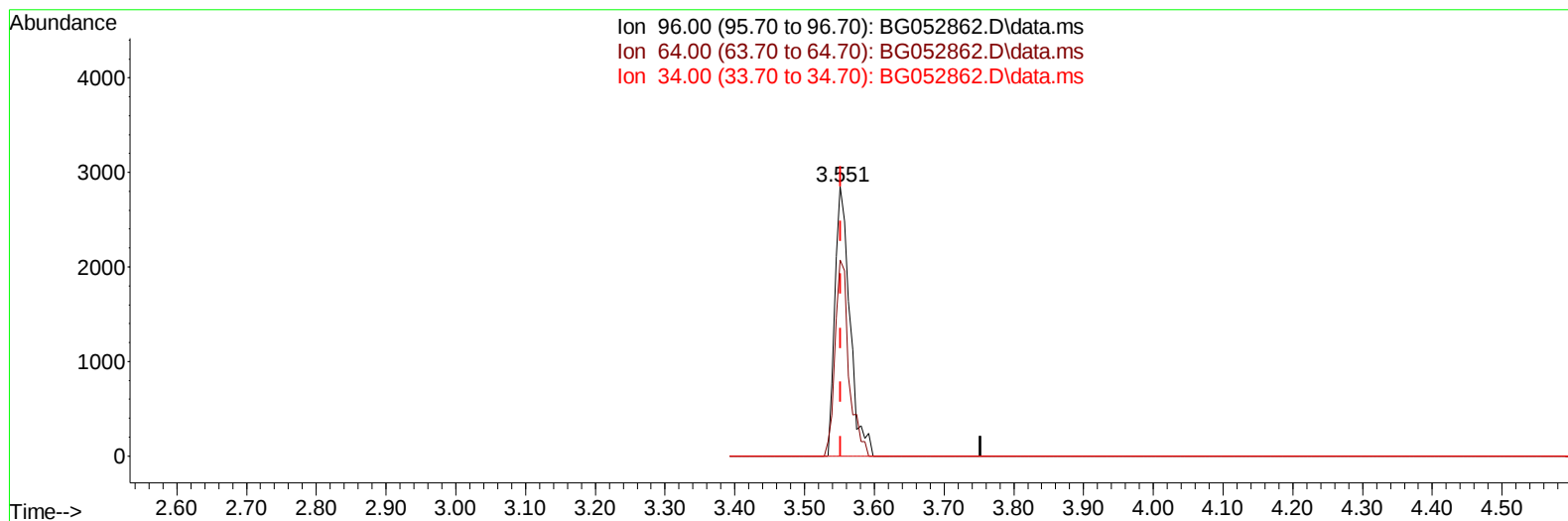
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TIC: BG052862.D\data.ms

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

3.551min (-0.001) 5.19 ng/uL m

response 4227

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	83.40	72.78
34.00	0.00	0.00
0.00	0.00	0.00

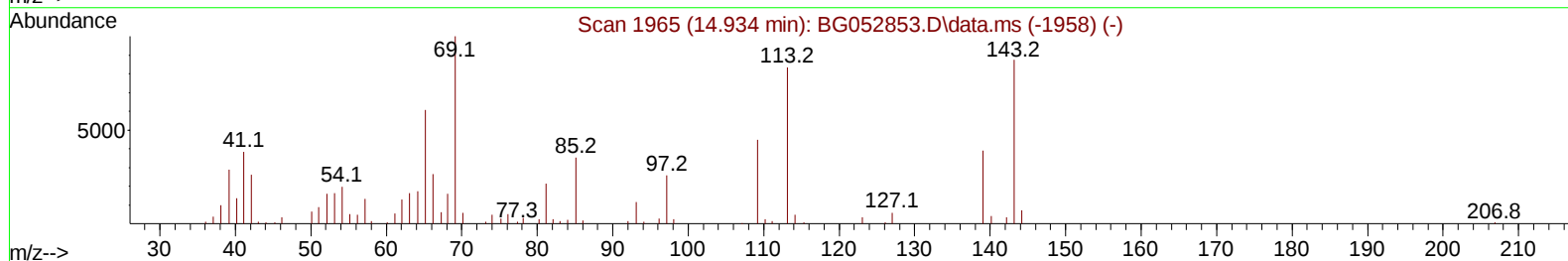
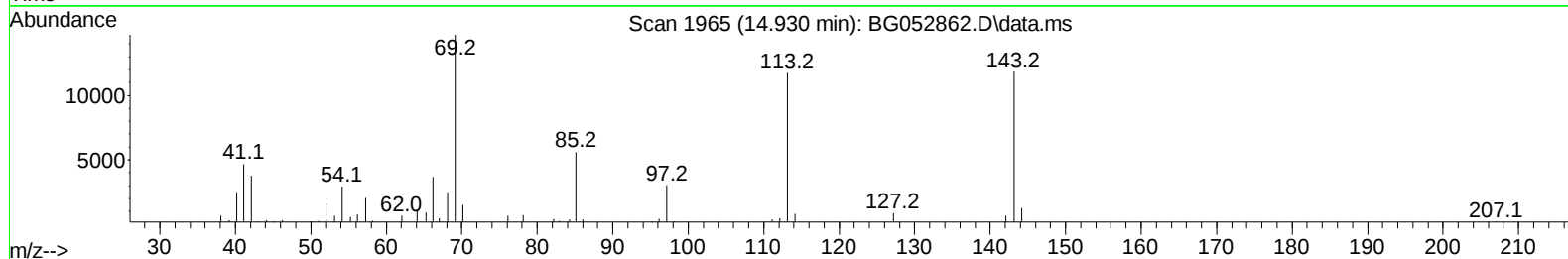
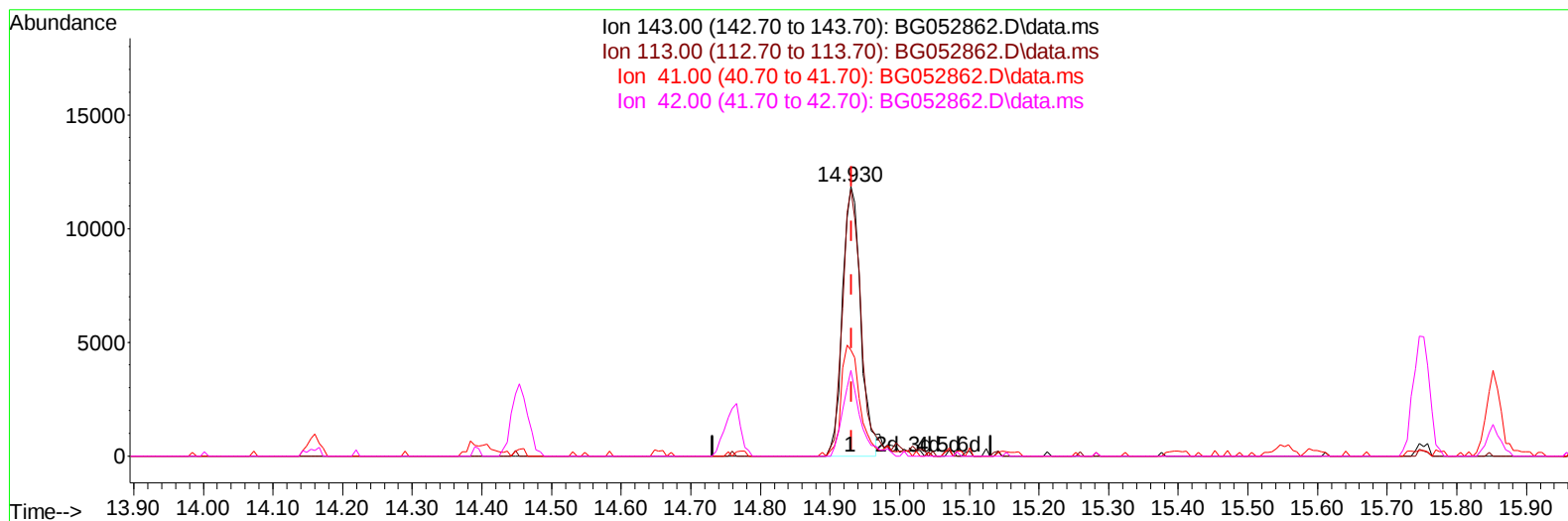
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TIC: BG052862.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.930min (-0.001) 15.49 ng/u1

response 21512

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	105.70	99.01
41.00	47.20	39.30
42.00	29.40	31.77

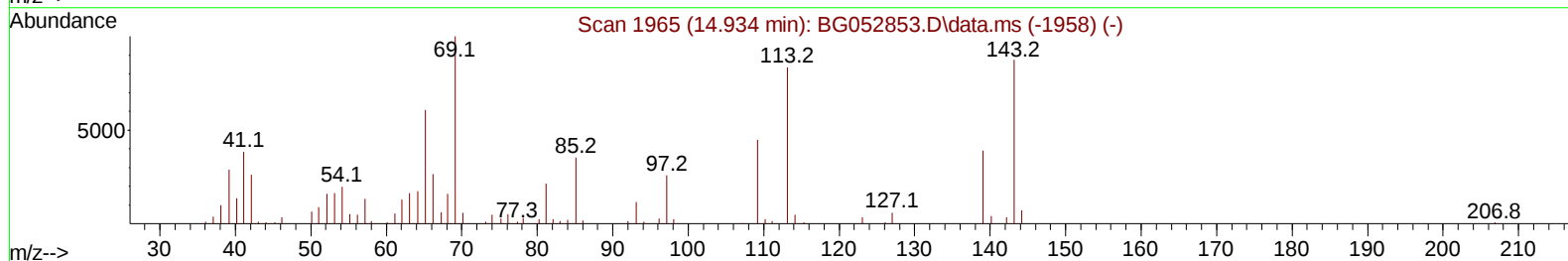
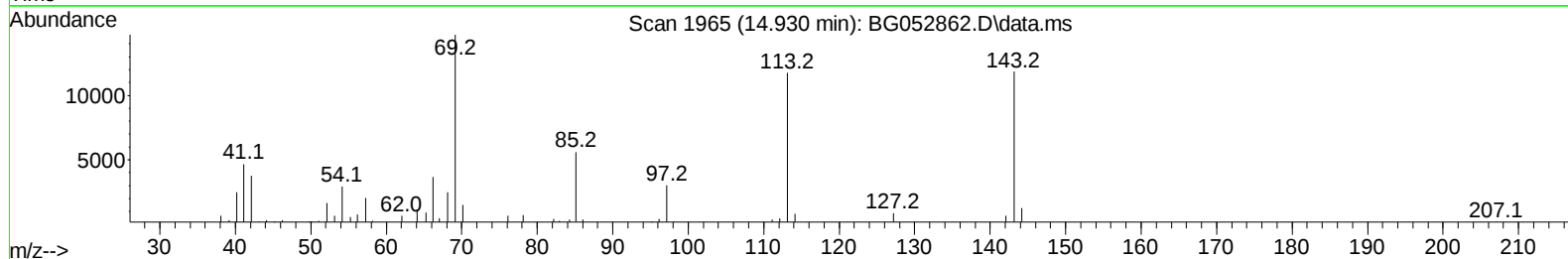
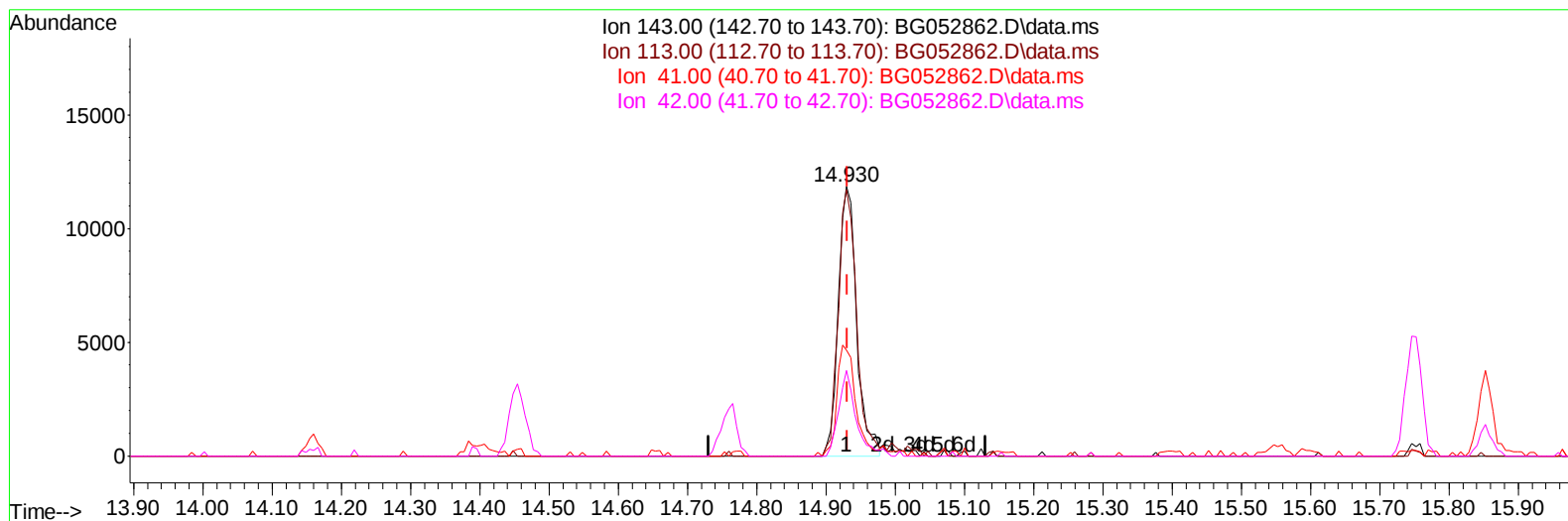
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Instrument :
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TIC: BG052862.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.930min (-0.001) 15.83 ng/ul m

response 21975

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	105.70	99.01
41.00	47.20	39.30
42.00	29.40	31.77

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.139	152	28794	20.000	ng/ul	0.00
20) Naphthalene-d8	10.953	136	131781	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.760	164	107494	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.503	188	268217	20.000	ng/ul	0.00
79) Chrysene-d12	21.791	240	299893	20.000	ng/ul	0.00
88) Perylene-d12	25.104	264	292561	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.551	96	4227m	5.185	ng/uL	0.00
4) Pyridine-d5	3.974	84	27447	13.487	ng/ul	0.00
7) Phenol-d5	7.287	99	36605	14.194	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.452	67	68166	43.277	ng/ul	0.00
11) 2-Chlorophenol-d4	7.669	132	59856	34.181	ng/ul	0.00
15) 4-Methylphenol-d8	8.832	113	49023	26.241	ng/ul	0.00
21) Nitrobenzene-d5	9.296	128	40869	43.083	ng/ul	0.00
24) 2-Nitrophenol-d4	10.025	143	47968	47.942	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.565	165	79986	36.747	ng/ul	0.00
31) 4-Chloroaniline-d4	11.076	131	102820	35.740	ng/ul	0.00
46) Dimethylphthalate-d6	14.160	166	355243	43.017	ng/ul	0.00
49) Acenaphthylene-d8	14.454	160	398357	39.697	ng/ul	0.00
54) 4-Nitrophenol-d4	14.930	143	21975m	15.827	ng/ul	0.00
60) Fluorene-d10	15.752	176	316431	43.639	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.852	200	67387	49.369	ng/ul	0.00
73) Anthracene-d10	17.603	188	578516	46.319	ng/ul	0.00
81) Pyrene-d10	19.888	212	708908	42.256	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.881	264	731175	48.369	ng/ul	0.00
Target Compounds						
78) Di-n-butylphthalate	18.455	149	18444	1.247	ng/ul	Qvalue 98

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

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