

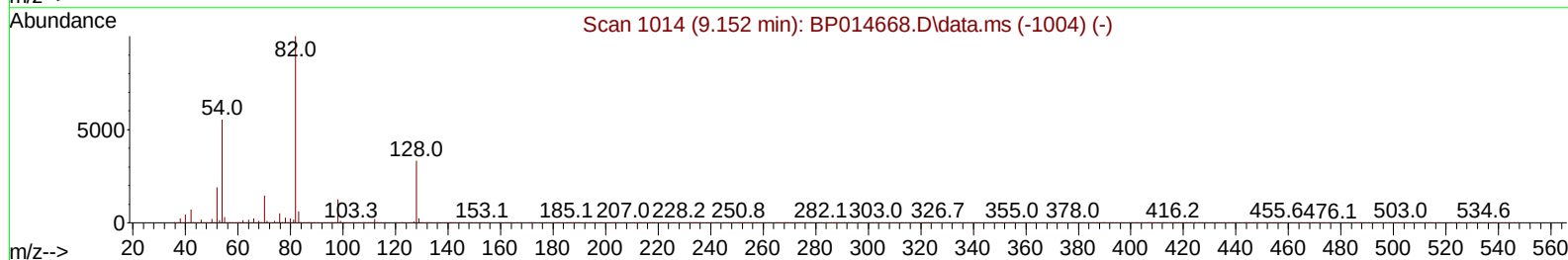
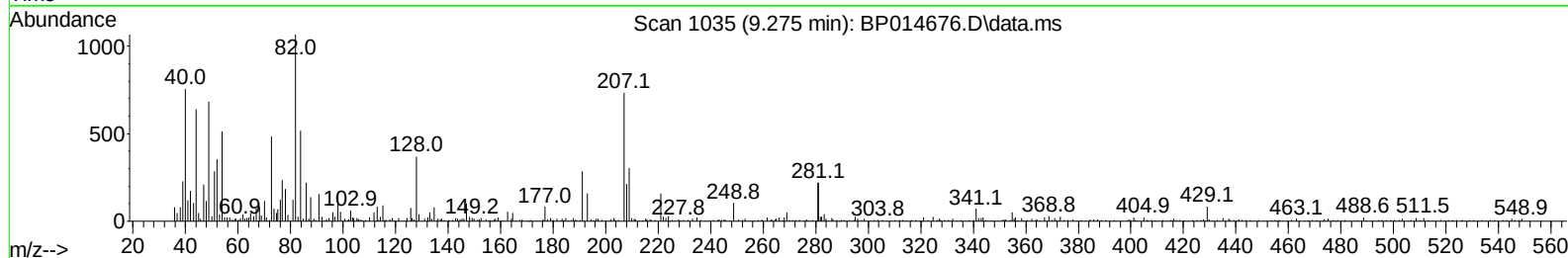
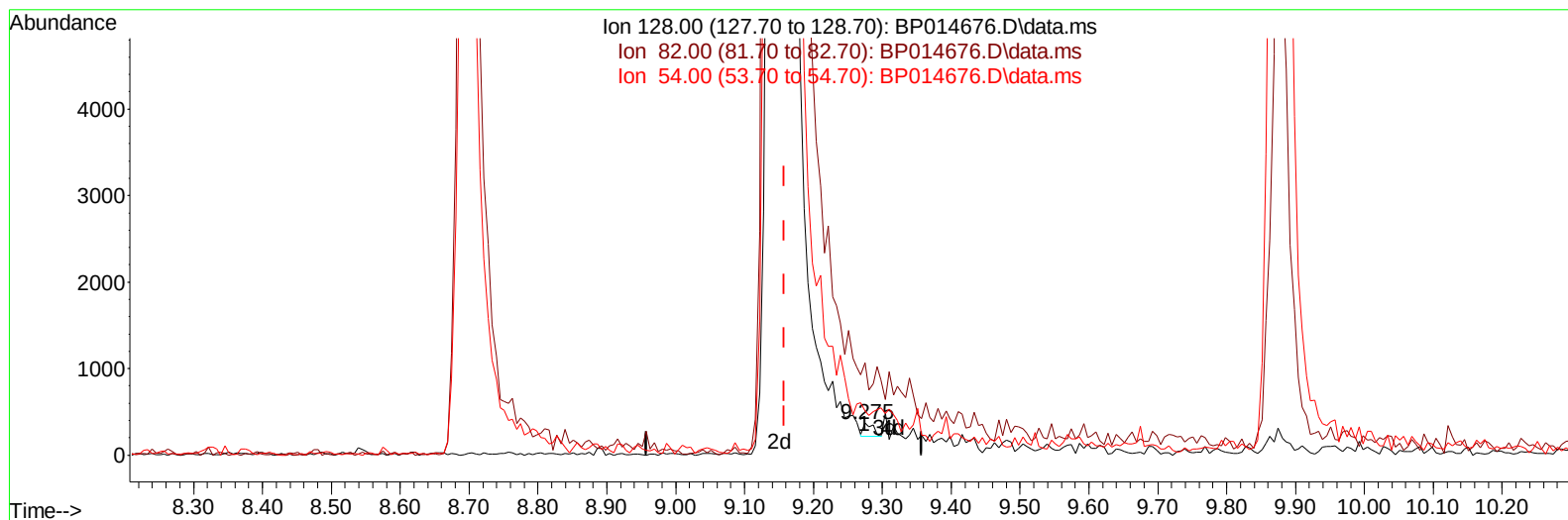
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041223\
Data File : BP014676.D
Acq On : 12 Apr 2023 16:39
Operator : CG/JU
Sample : 02251-17
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
COMD5

Manual IntegrationsAPPROVED

Quant Time: Apr 13 00:50:01 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP032823.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue Apr 11 01:21:43 2023
Response via : Initial Calibration

Reviewed By :Christian Giraldo 04/13/2023
Supervised By :Jagrut Upadhyay 04/13/2023



TIC: BP014676.D\data.ms

(21) Nitrobenzene-d5 (S)

9.275min (+ 0.117) 0.04 ng/uL

response 144

Ion	Exp%	Act%
128.00	100.00	100.00
82.00	278.40	288.11
54.00	139.60	138.38
0.00	0.00	0.00

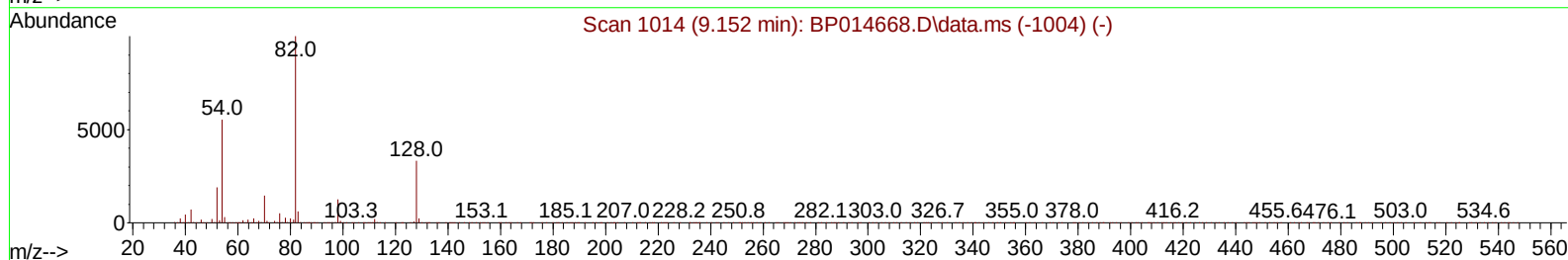
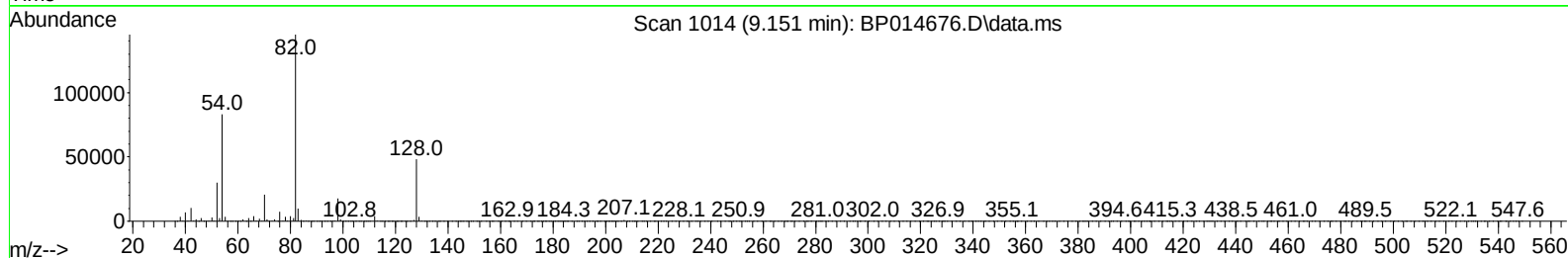
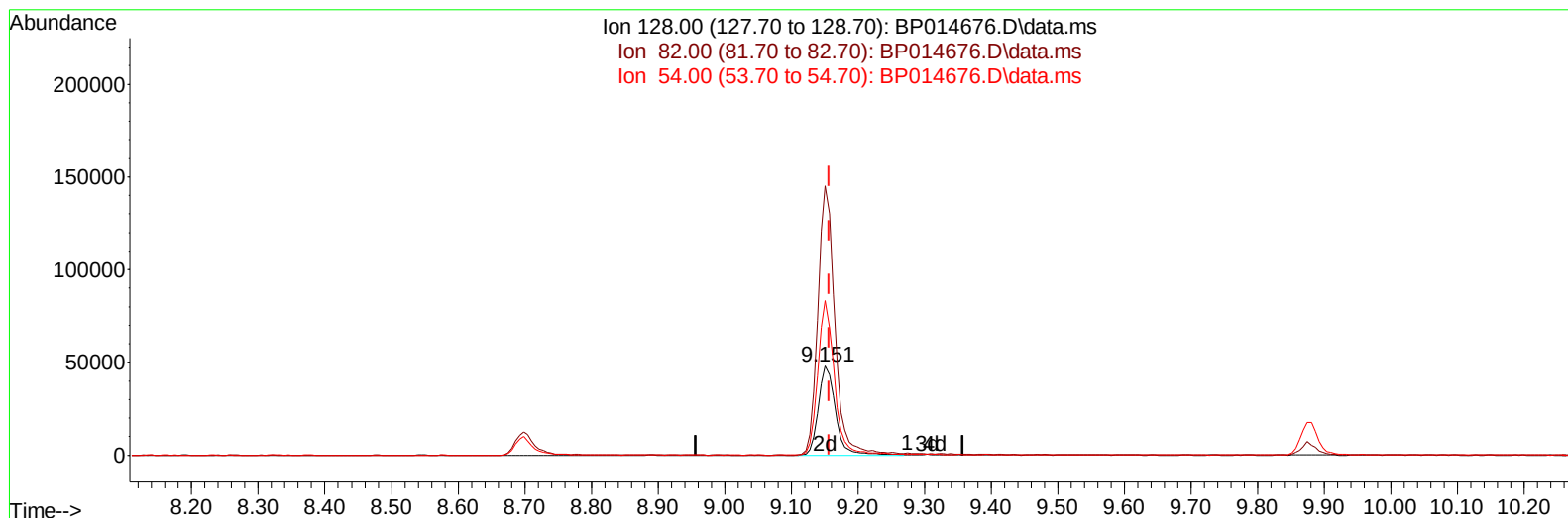
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(21) Nitrobenzene-d5 (S)

9.151min (-0.006) 25.82 ng/u1 m

response 85965

Ion	Exp%	Act%
128.00	100.00	100.00
82.00	278.40	302.33
54.00	139.60	173.53#
0.00	0.00	0.00

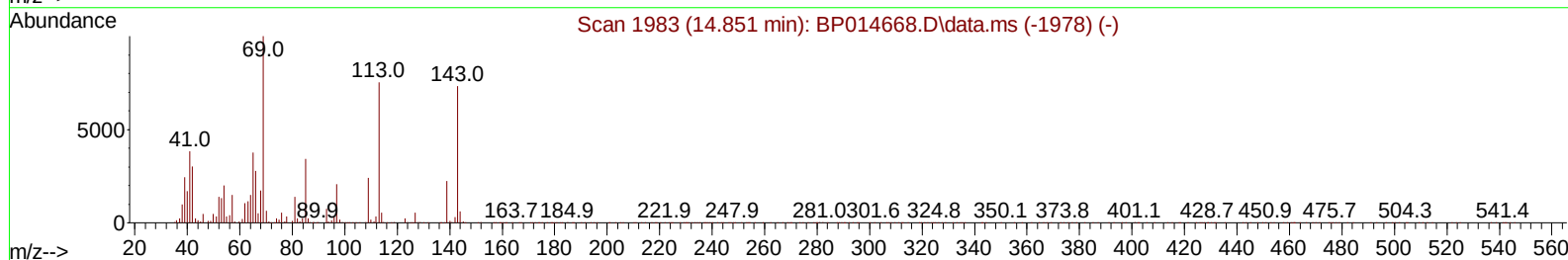
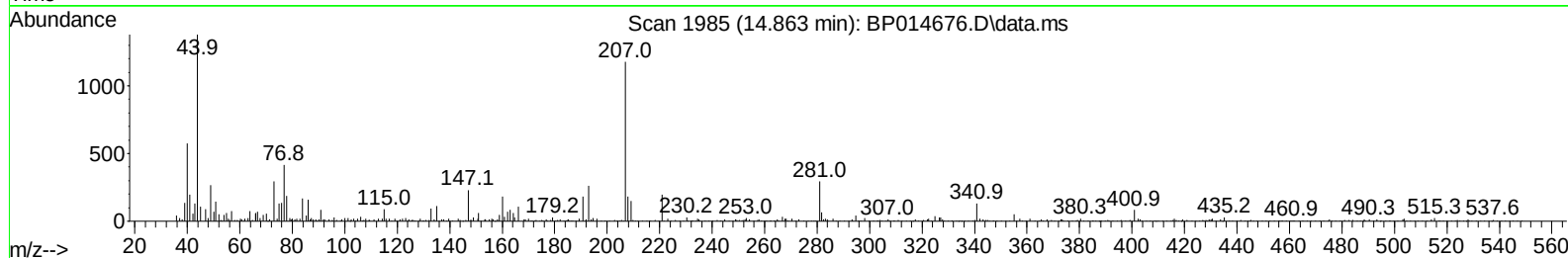
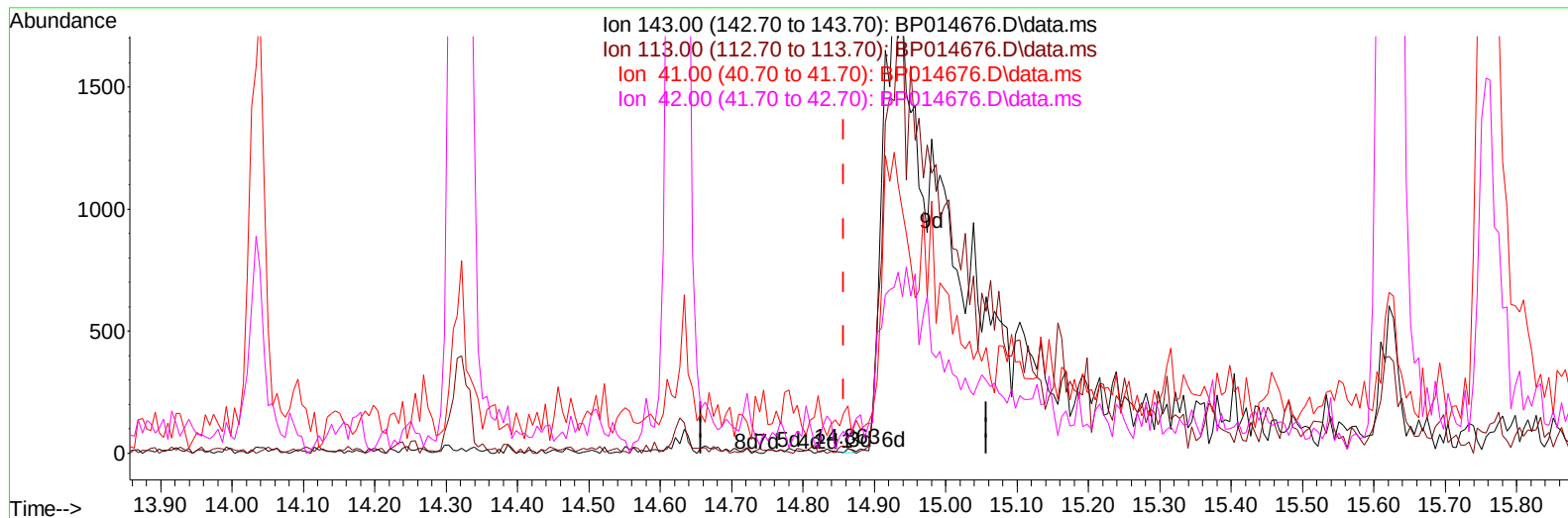
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Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
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TIC: BP014676.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.863min (+ 0.006) 0.00 ng/ul

response 10

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	96.80	116.67#
41.00	44.20	1083.33#
42.00	31.50	311.11#

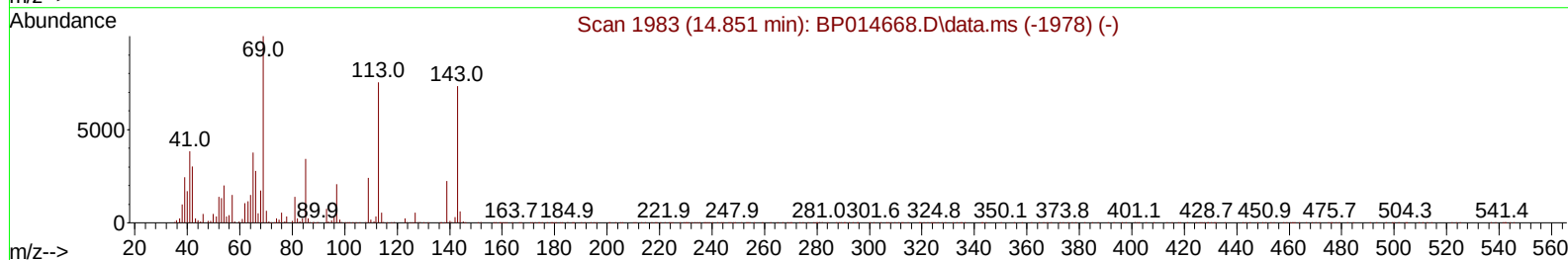
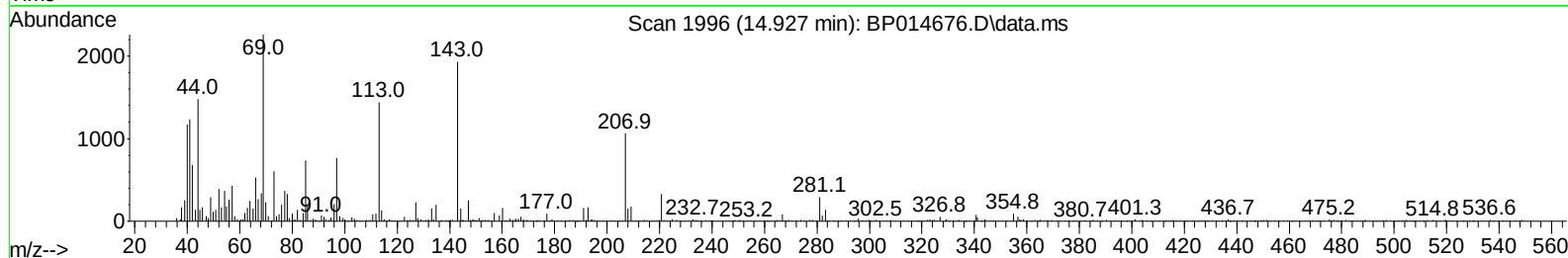
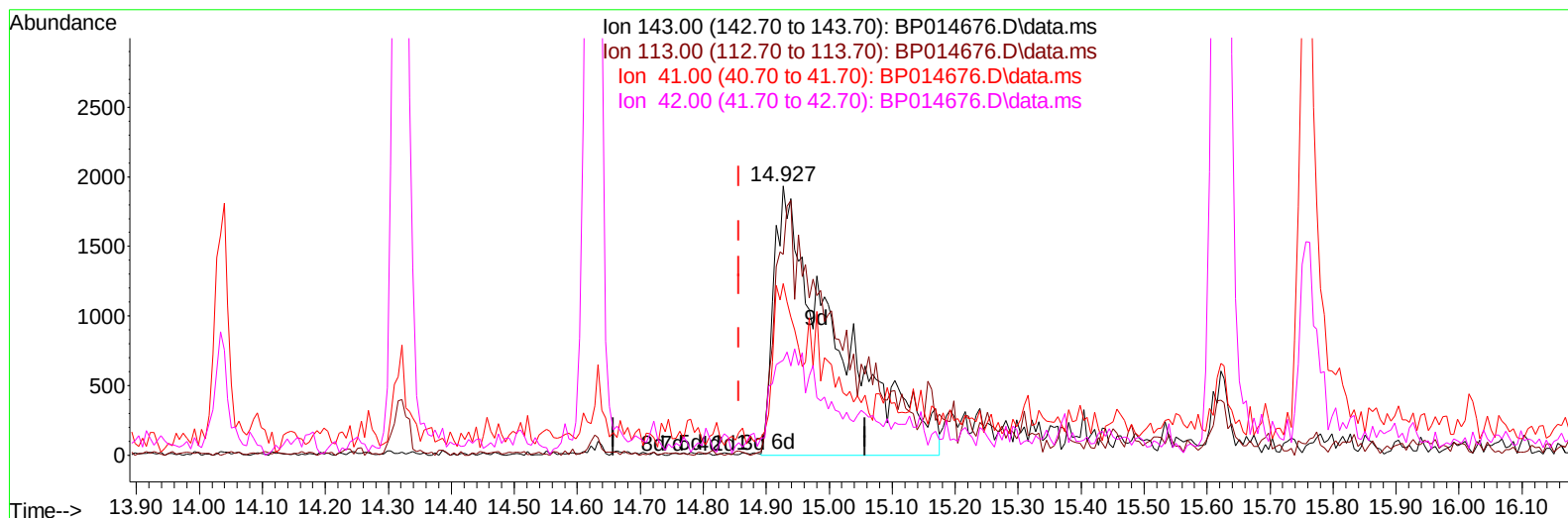
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(54) 4-Nitrophenol-d4 (S)

14.927min (+ 0.070) 4.20 ng/ul m

response 13163

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	96.80	74.70#
41.00	44.20	63.84#
42.00	31.50	35.18

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.975	152	99146	20.000	ng/ul	0.00
20) Naphthalene-d8	10.792	136	412468	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.627	164	255567	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.392	188	564691	20.000	ng/ul	0.00
79) Chrysene-d12	21.480	240	572874	20.000	ng/ul	0.00
88) Perylene-d12	23.980	264	672480	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.363	96	14367	5.619	ng/uL	0.00
4) Pyridine-d5	3.810	84	40272	6.147	ng/ul	0.02
7) Phenol-d5	7.145	99	57629	6.303	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.298	67	172115	29.338	ng/ul	0.00
11) 2-Chlorophenol-d4	7.504	132	143901	21.774	ng/ul	0.00
15) 4-Methylphenol-d8	8.698	113	101247	13.727	ng/ul	0.00
21) Nitrobenzene-d5	9.151	128	85965m	25.817	ng/ul	0.00
24) 2-Nitrophenol-d4	9.875	143	92201	24.182	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.422	165	151532	21.822	ng/ul	0.00
31) 4-Chloroaniline-d4	10.945	131	188432	20.601	ng/ul	0.00
46) Dimethylphthalate-d6	14.033	166	549305	26.508	ng/ul	0.00
49) Acenaphthylene-d8	14.322	160	610925	26.570	ng/ul	0.00
54) 4-Nitrophenol-d4	14.927	143	13163m	4.205	ng/ul	0.07
60) Fluorene-d10	15.621	176	472839	27.363	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.757	200	73157	17.607	ng/ul	0.00
73) Anthracene-d10	17.492	188	758711	27.763	ng/ul	0.00
81) Pyrene-d10	19.721	212	947907	24.634	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.821	264	1012148	28.329	ng/ul	0.00

Target Compounds Qvalue

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

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