

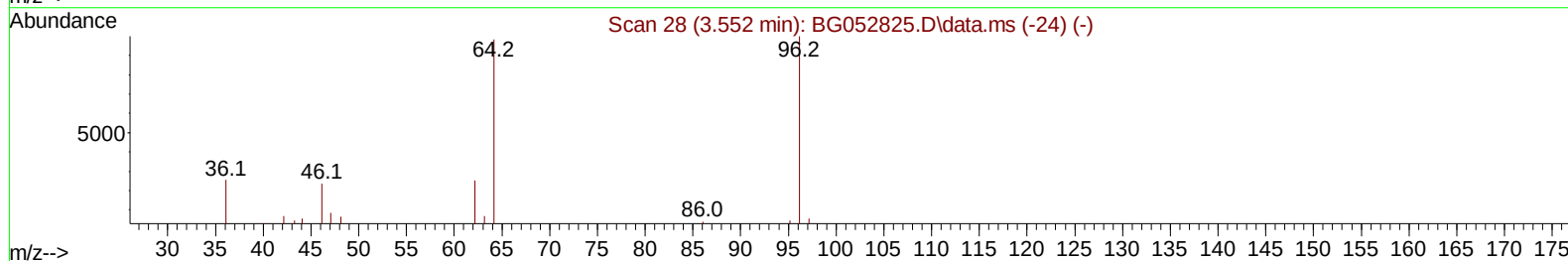
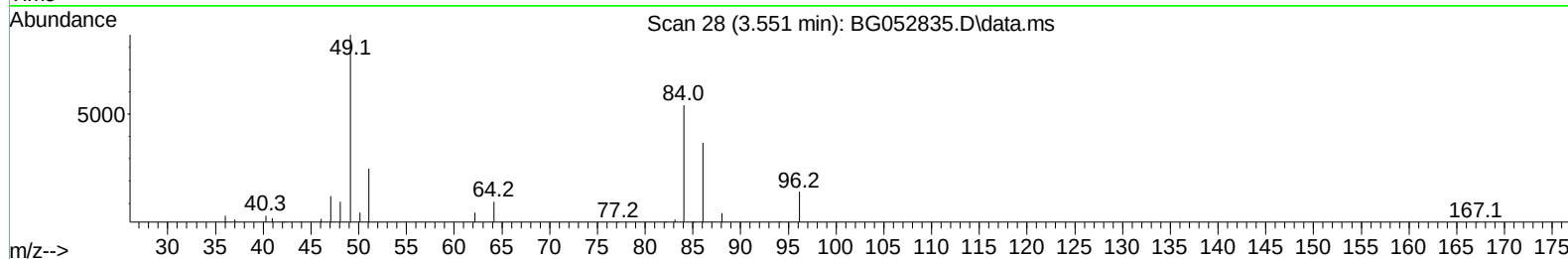
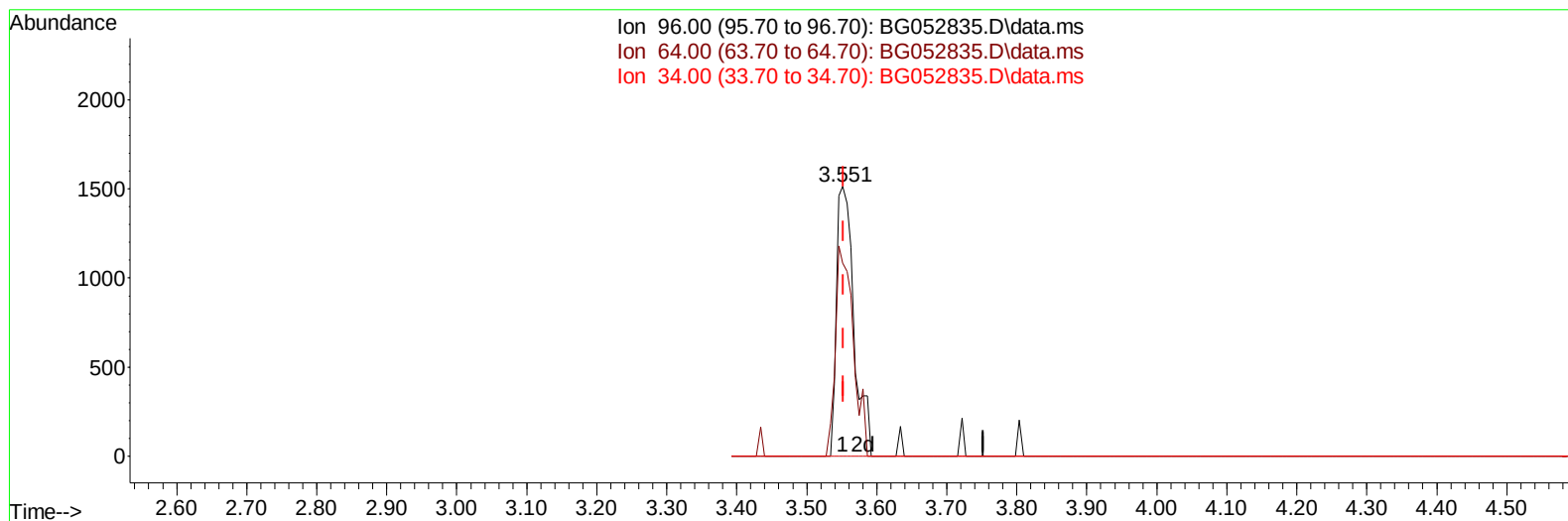
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG032522\
Data File : BG052835.D
Acq On : 25 Mar 2022 18:40
Operator : CG/JU
Sample : N2079-12
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
EW5Z5

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 03/28/2022
Supervised By :mohammad ahmed 03/29/2022

Quant Time: Mar 26 00:46:28 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
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TIC: BG052835.D\data.ms

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

3.551min (-0.001) 3.95 ng/uL m

response 2623

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	83.40	71.60
34.00	0.00	0.00
0.00	0.00	0.00

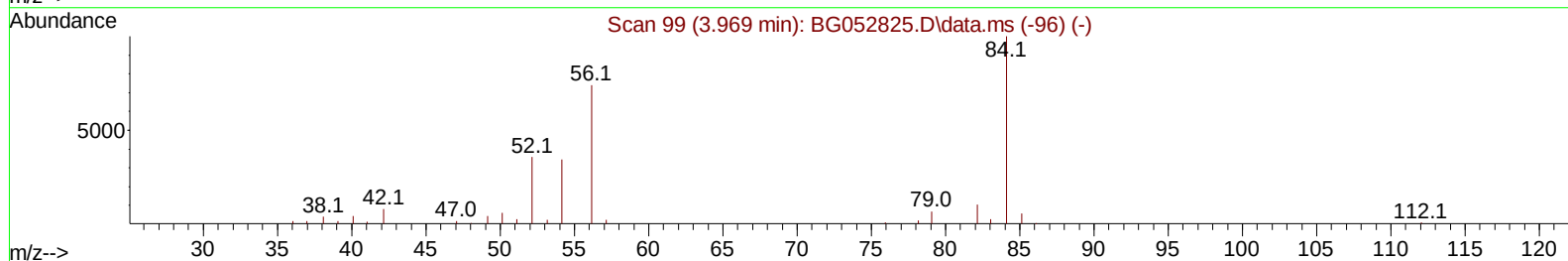
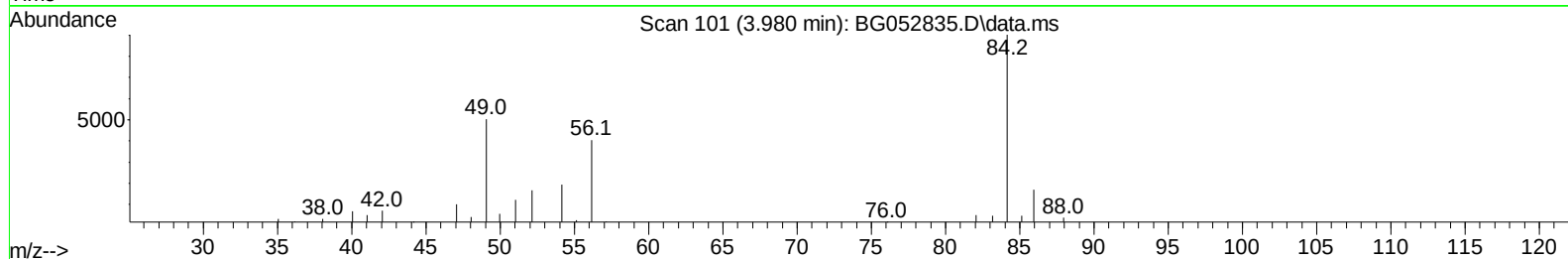
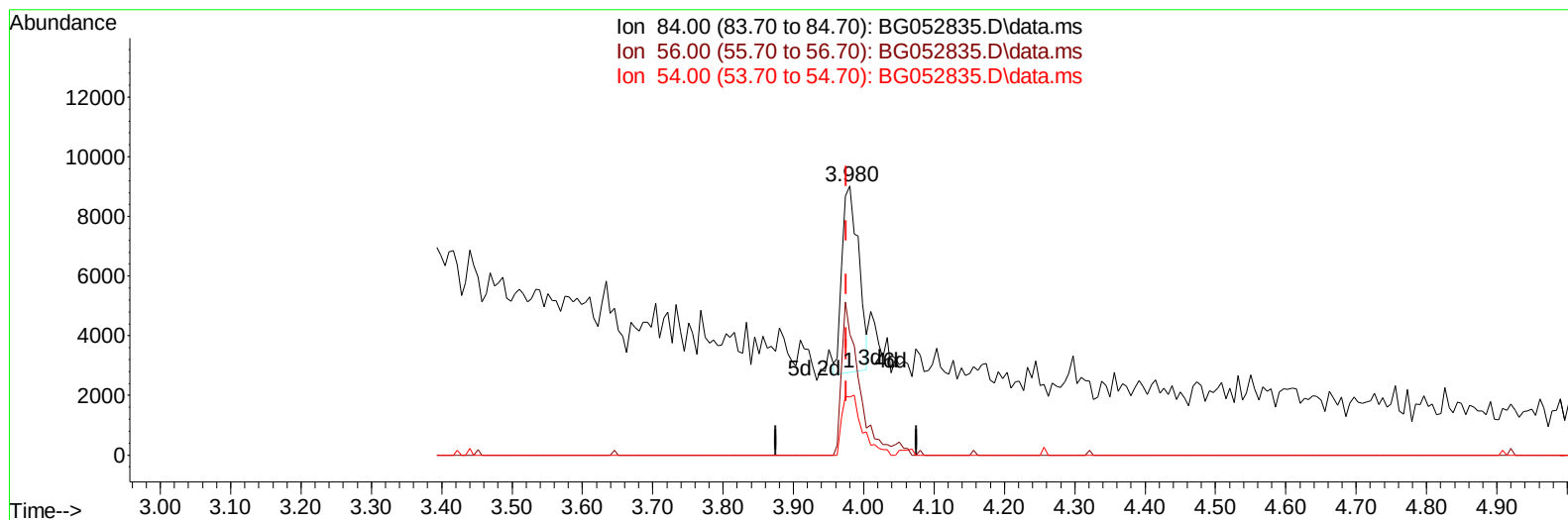
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(4) Pyridine-d5 (S)

3.980min (+ 0.005) 6.16 ng/ul

response 10213

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	65.20	44.89#
54.00	28.70	21.55#
0.00	0.00	0.00

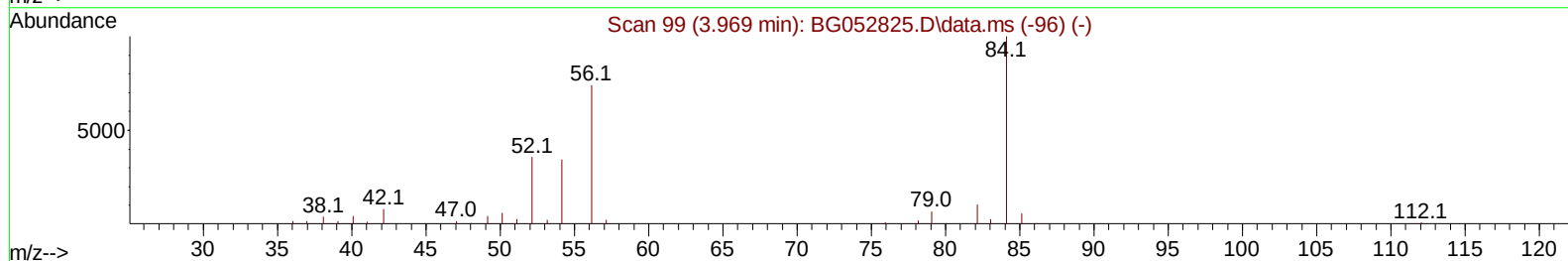
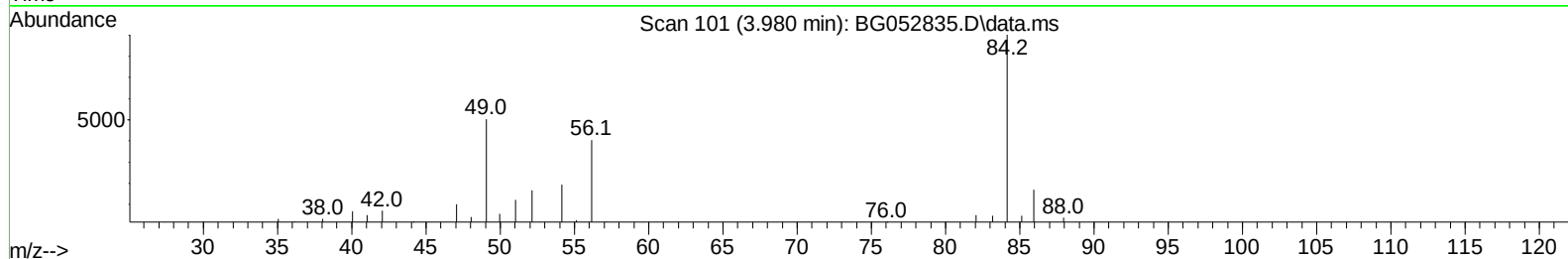
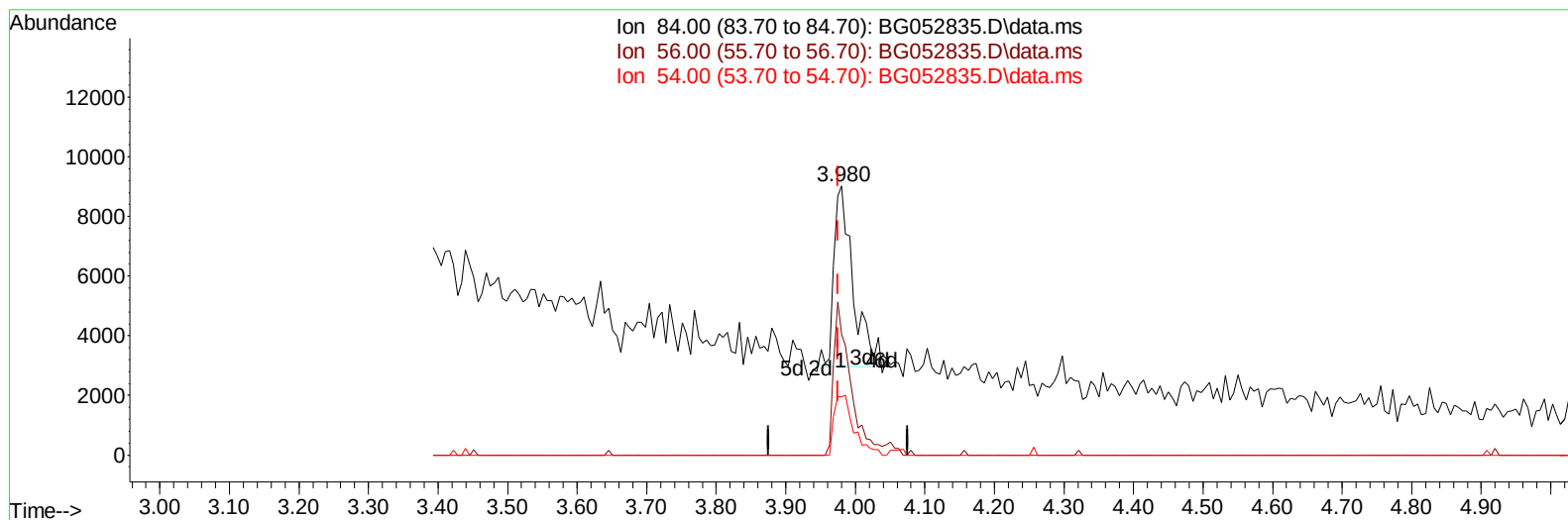
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TIC: BG052835.D\data.ms

(4) Pyridine-d5 (S)

3.980min (+ 0.005) 6.75 ng/ul m

response 11186

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	65.20	44.89#
54.00	28.70	21.55#
0.00	0.00	0.00

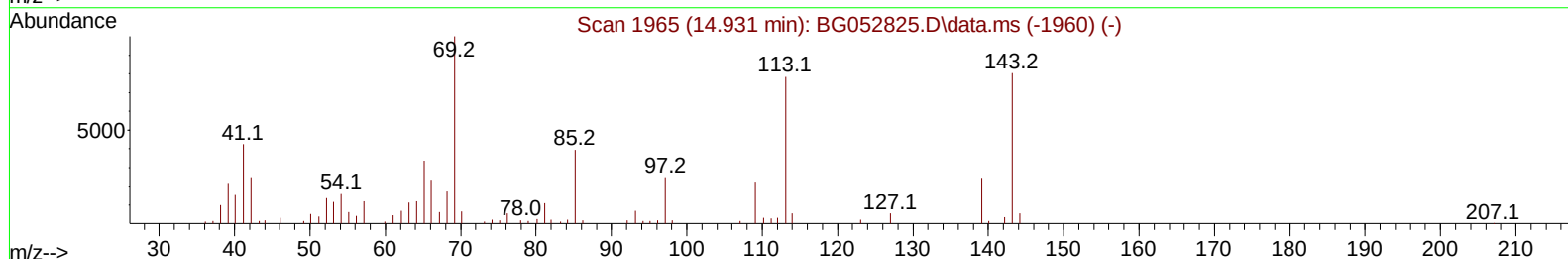
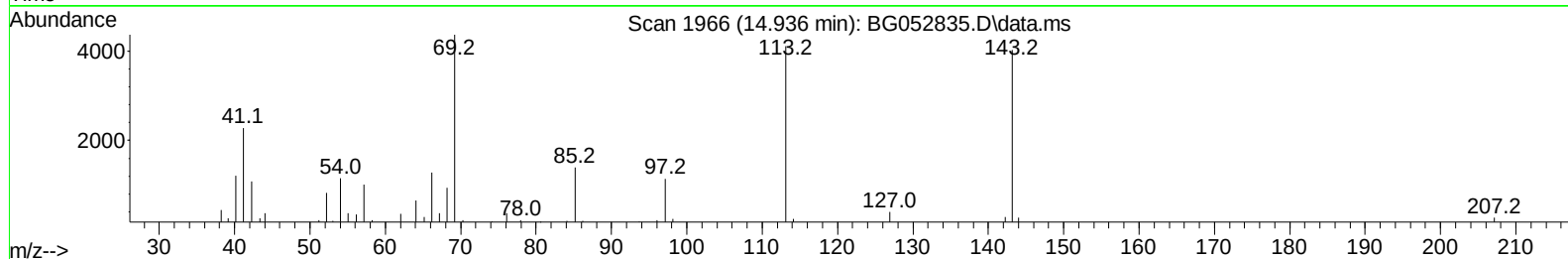
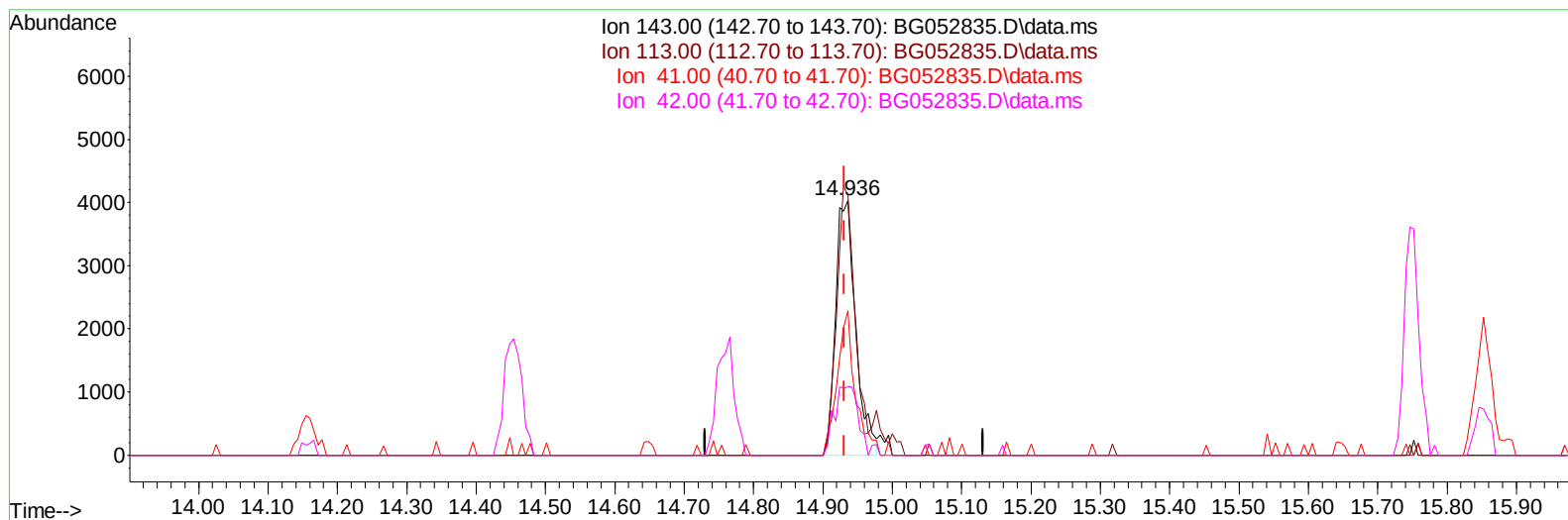
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TIC: BG052835.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.936min (+ 0.005) 7.11 ng/u1

response 8140

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	105.70	102.03
41.00	47.20	56.58
42.00	29.40	26.75

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.145	152	23458	20.000	ng/ul	0.00
20) Naphthalene-d8	10.953	136	108143	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.760	164	88690	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.503	188	224711	20.000	ng/ul	0.00
79) Chrysene-d12	21.791	240	255380	20.000	ng/ul	0.00
88) Perylene-d12	25.104	264	259522	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.551	96	2623m	3.950	ng/uL	0.00
4) Pyridine-d5	3.980	84	11186m	6.747	ng/ul	0.00
7) Phenol-d5	7.287	99	13790	6.563	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.458	67	45818	35.706	ng/ul	0.00
11) 2-Chlorophenol-d4	7.669	132	35071	24.583	ng/ul	0.00
15) 4-Methylphenol-d8	8.838	113	26138	17.174	ng/ul	0.00
21) Nitrobenzene-d5	9.296	128	25283	32.478	ng/ul	0.00
24) 2-Nitrophenol-d4	10.025	143	29637	36.095	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.565	165	49697	27.822	ng/ul	0.00
31) 4-Chloroaniline-d4	11.076	131	59272	25.106	ng/ul	0.00
46) Dimethylphthalate-d6	14.160	166	224480	32.946	ng/ul	0.00
49) Acenaphthylene-d8	14.454	160	259914	31.393	ng/ul	0.00
54) 4-Nitrophenol-d4	14.936	143	8319m	7.262	ng/ul	0.00
60) Fluorene-d10	15.752	176	204528	34.187	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.852	200	41892	36.633	ng/ul	0.00
73) Anthracene-d10	17.603	188	388937	37.169	ng/ul	0.00
81) Pyrene-d10	19.888	212	498091	34.865	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.887	264	494973	36.912	ng/ul	0.01

Target Compounds Qvalue

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

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