

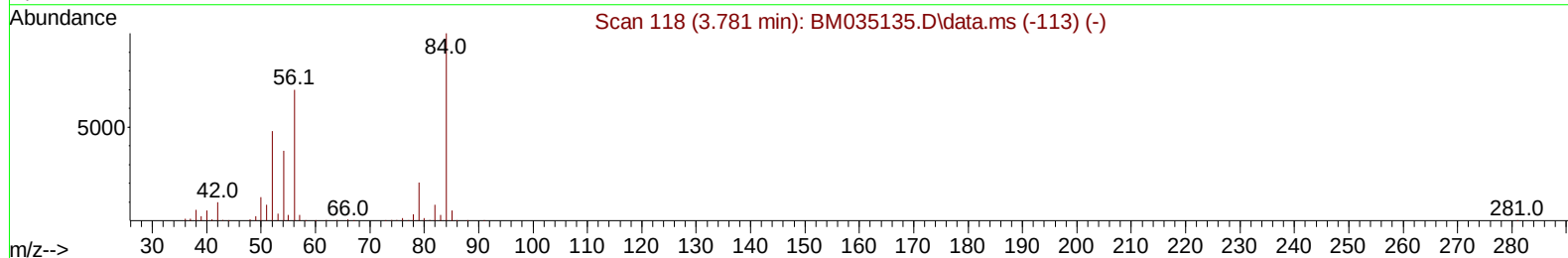
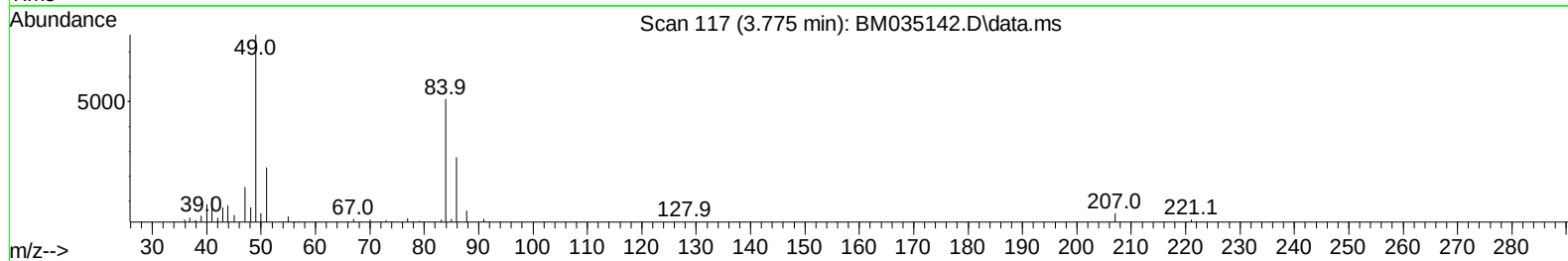
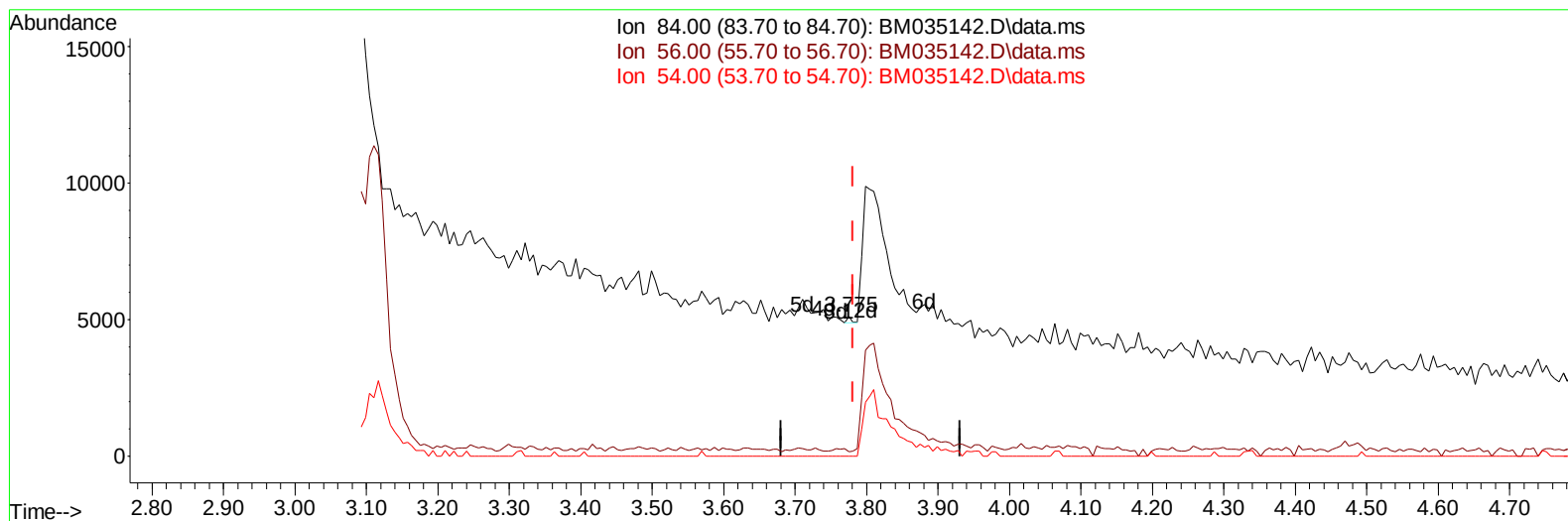
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM051822\
Data File : BM035142.D
Acq On : 18 May 2022 13:55
Operator : CG/JU
Sample : N2766-20
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
EW4H0

Manual IntegrationsAPPROVED

Quant Time: May 19 00:49:38 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM051722.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue May 17 16:13:52 2022
Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 05/19/2022
Supervised By :mohammad ahmed 05/19/2022



TIC: BM035142.D\data.ms

(4) Pyridine-d5 (S)

3.775min (-0.006) 0.01 ng/ul

response 114

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	68.20	3.21#
54.00	34.50	0.00#
0.00	0.00	0.00

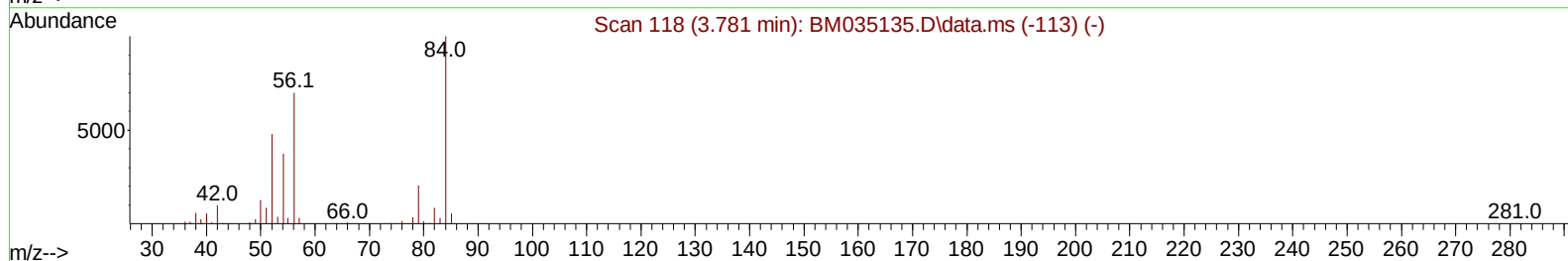
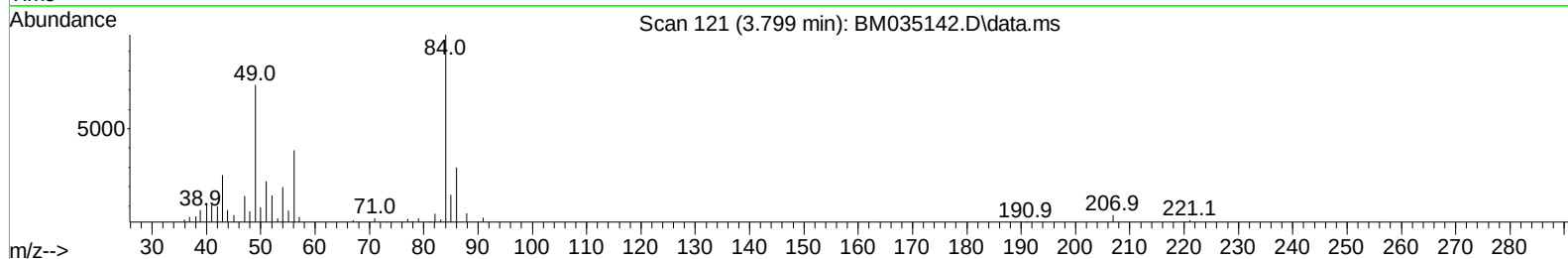
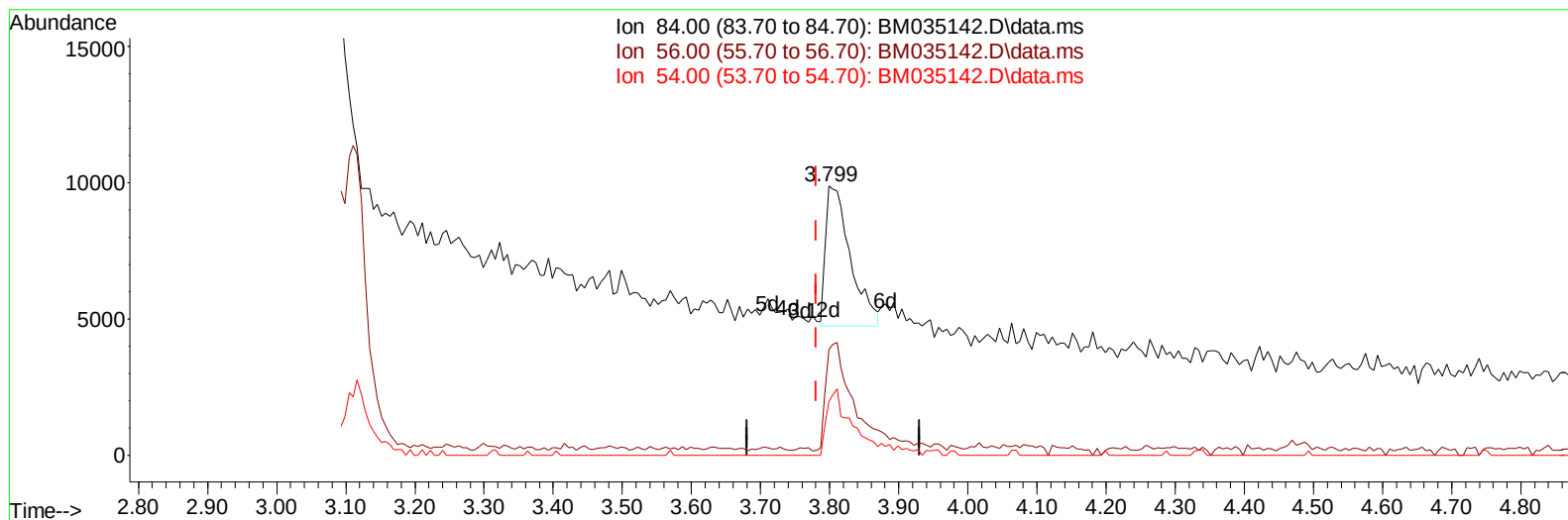
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(4) Pyridine-d5 (S)

3.799min (+ 0.018) 1.16 ng/ul m

response 12675

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	68.20	39.35#
54.00	34.50	19.84#
0.00	0.00	0.00

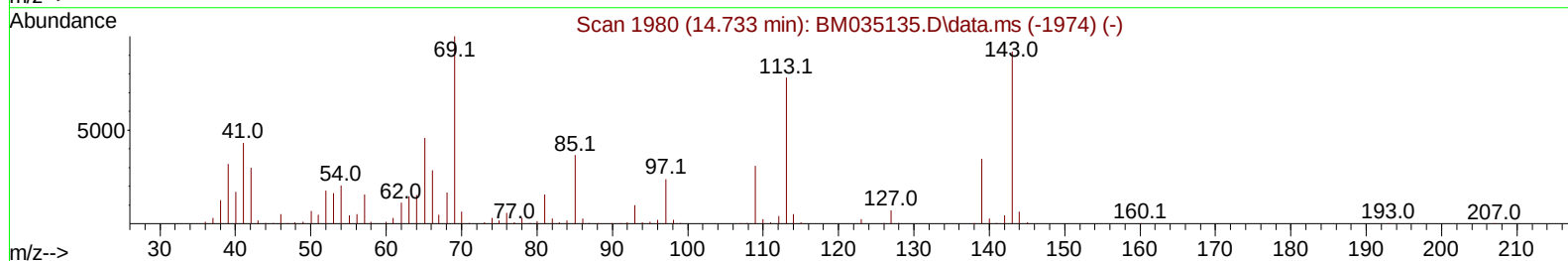
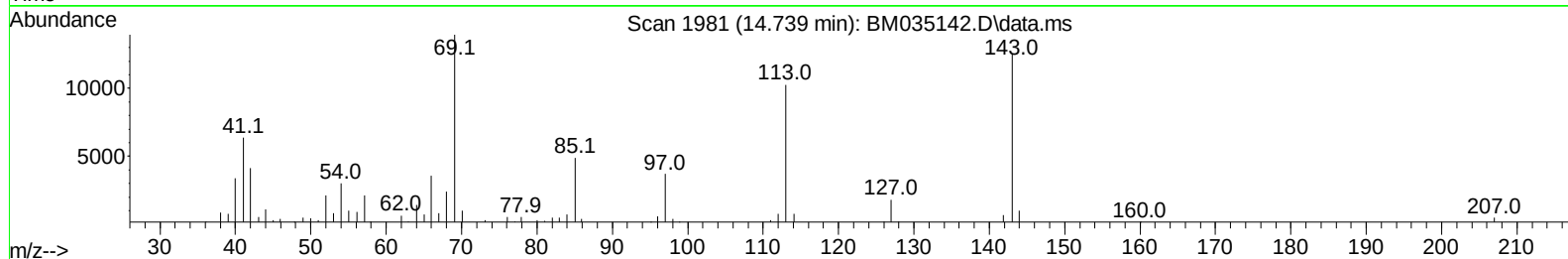
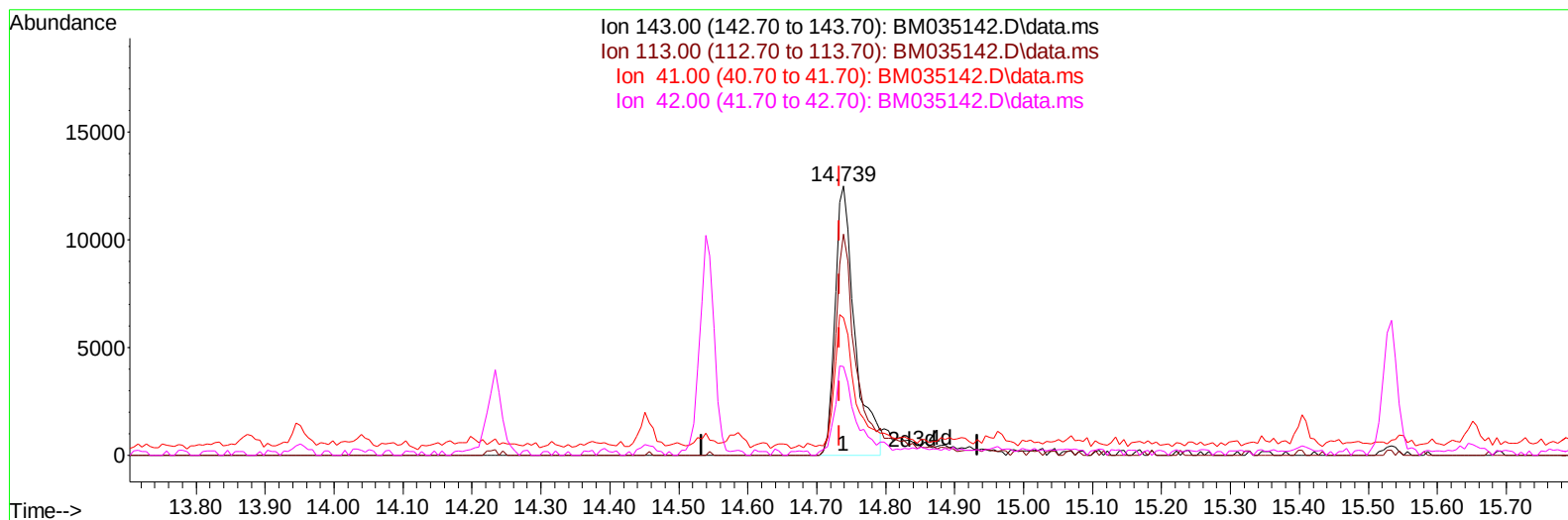
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TIC: BM035142.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.739min (+ 0.006) 4.97 ng/u1

response 25029

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	80.00	82.16
41.00	42.20	51.08#
42.00	28.30	32.92

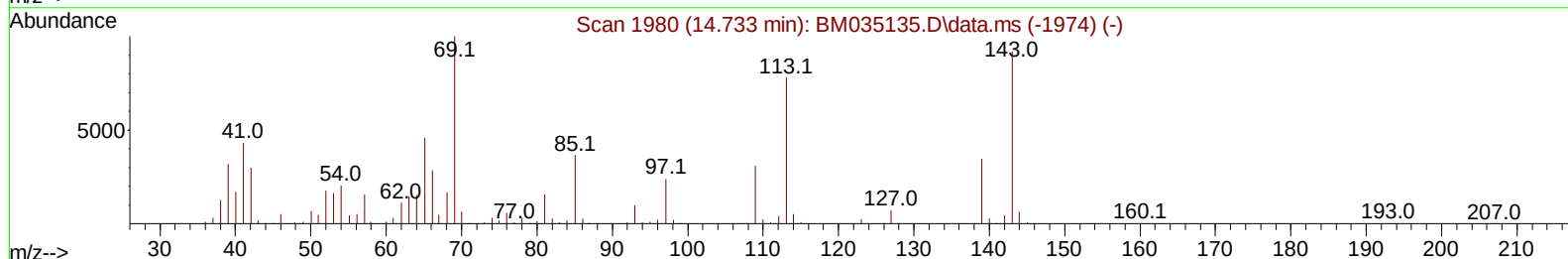
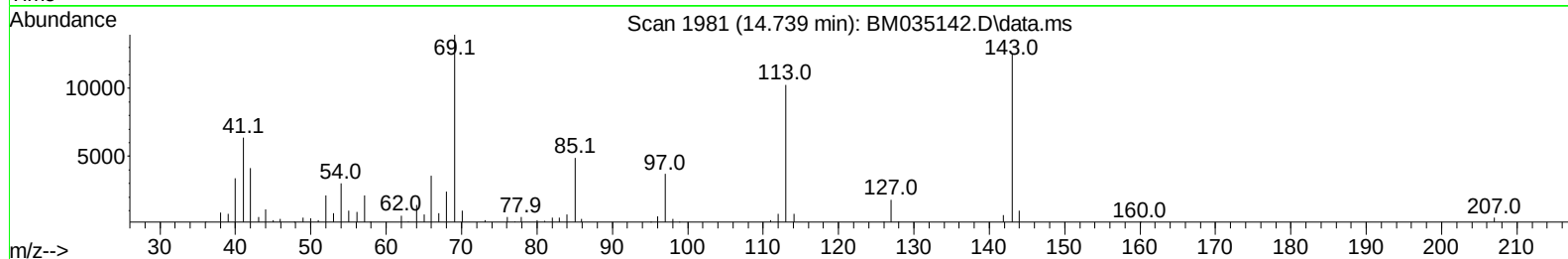
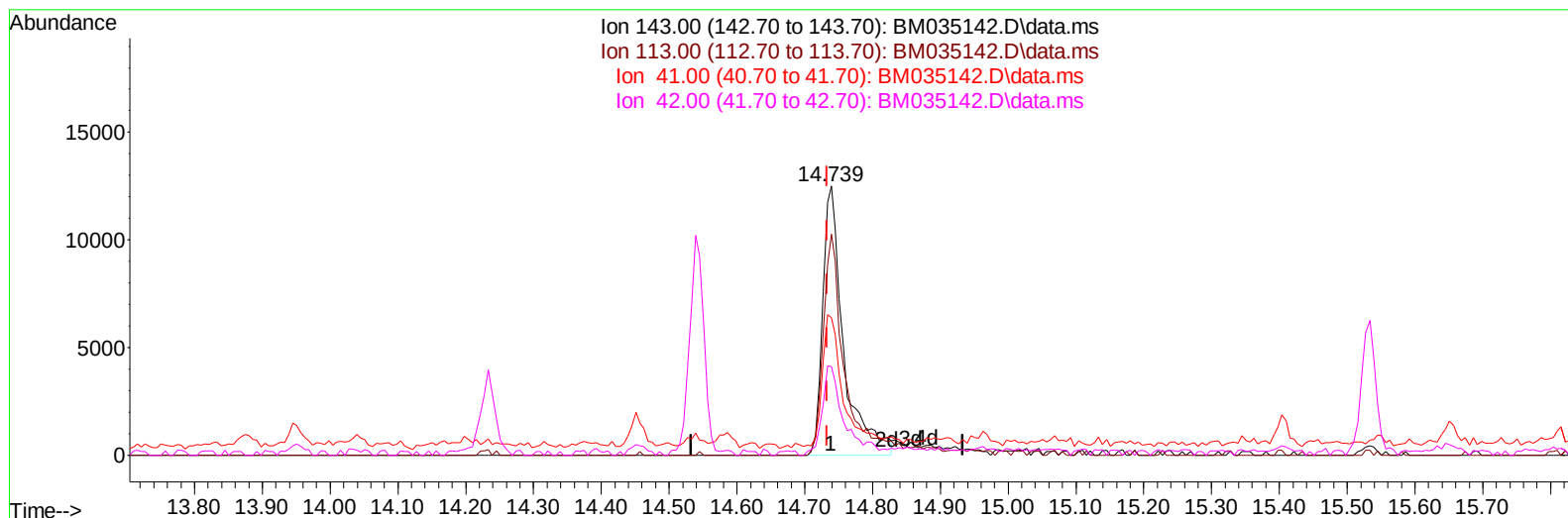
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Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
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TIC: BM035142.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.739min (+ 0.006) 5.34 ng/ul m

response 26905

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	80.00	82.16
41.00	42.20	51.08#
42.00	28.30	32.92

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.910	152	174321	20.000	ng/ul	0.00
20) Naphthalene-d8	10.710	136	670230	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.539	164	389747	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.286	188	763696	20.000	ng/ul	0.00
79) Chrysene-d12	21.468	240	702990	20.000	ng/ul	0.00
88) Perylene-d12	23.856	264	737336	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.375	96	7994	1.895	ng/uL	0.01
4) Pyridine-d5	3.799	84	12675m	1.165	ng/ul	0.02
7) Phenol-d5	7.075	99	132154	10.287	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.240	67	90800	11.202	ng/ul	0.00
11) 2-Chlorophenol-d4	7.440	132	123688	10.860	ng/ul	0.00
15) 4-Methylphenol-d8	8.616	113	101023	9.361	ng/ul	0.00
21) Nitrobenzene-d5	9.063	128	55567	10.574	ng/ul	0.00
24) 2-Nitrophenol-d4	9.786	143	59241	10.069	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.328	165	105240	10.056	ng/ul	0.00
31) 4-Chloroaniline-d4	10.845	131	113168	7.842	ng/ul	0.00
46) Dimethylphthalate-d6	13.951	166	325736	11.471	ng/ul	0.00
49) Acenaphthylene-d8	14.233	160	380814	11.221	ng/ul	0.00
54) 4-Nitrophenol-d4	14.739	143	26905m	5.344	ng/ul	0.00
60) Fluorene-d10	15.533	176	281591	11.722	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.651	200	13177	2.672	ng/ul	0.00
73) Anthracene-d10	17.386	188	406759	11.443	ng/ul	0.00
81) Pyrene-d10	19.674	212	455056	12.154	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.703	264	389379	10.420	ng/ul	0.00
Target Compounds						
					Qvalue	
16) Acetophenone	8.734	105	26225	1.569	ng/ul	96

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

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