

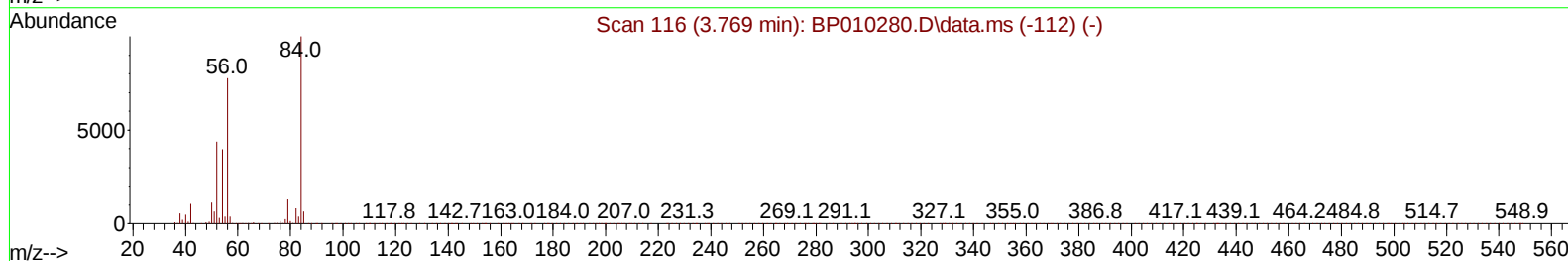
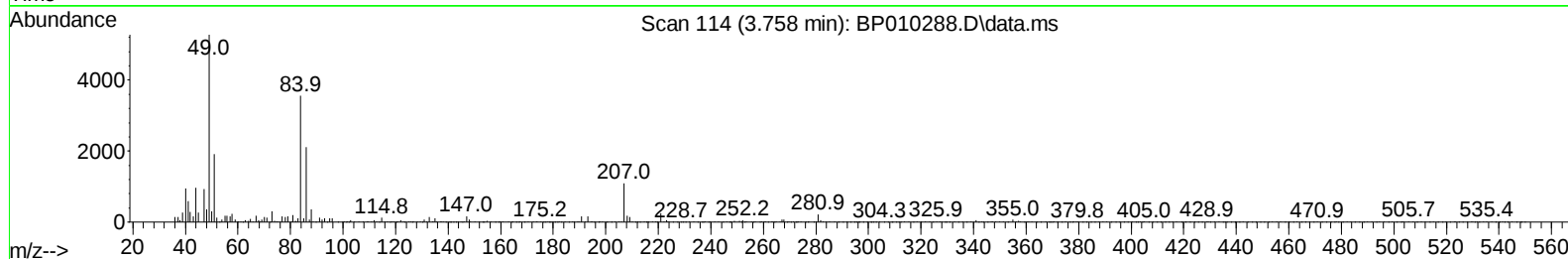
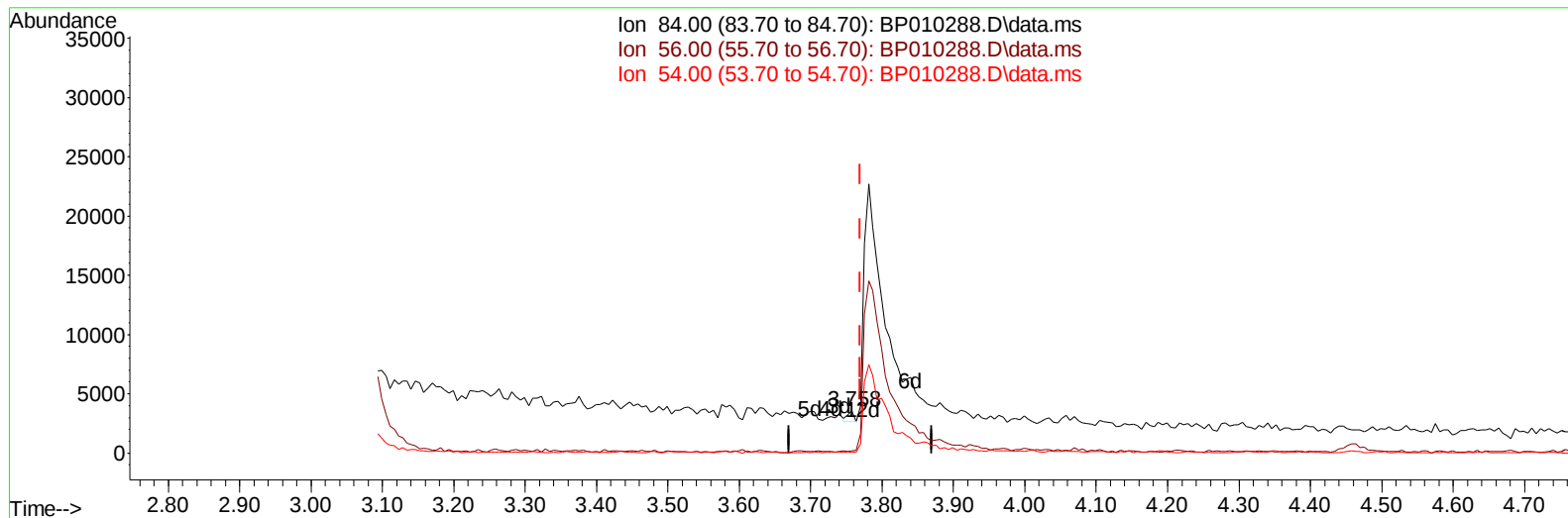
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051122\
 Data File : BP010288.D
 Acq On : 11 May 2022 14:16
 Operator : CG/JU
 Sample : N2632-01
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 EW4A3

Manual IntegrationsAPPROVED

Quant Time: May 12 01:13:16 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP051022.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu May 12 01:10:15 2022
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 05/12/2022
 Supervised By :Yogesh Patel 05/17/2022



TIC: BP010288.D\data.ms

(4) Pyridine-d5 (S)

3.758min (-0.012) 0.07 ng/u1

response 543

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	30.50	4.84#
54.00	24.80	2.17#
0.00	0.00	0.00

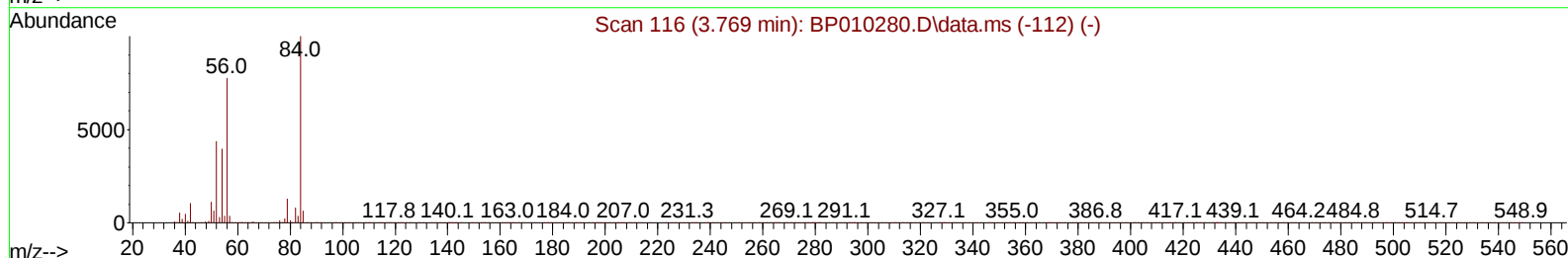
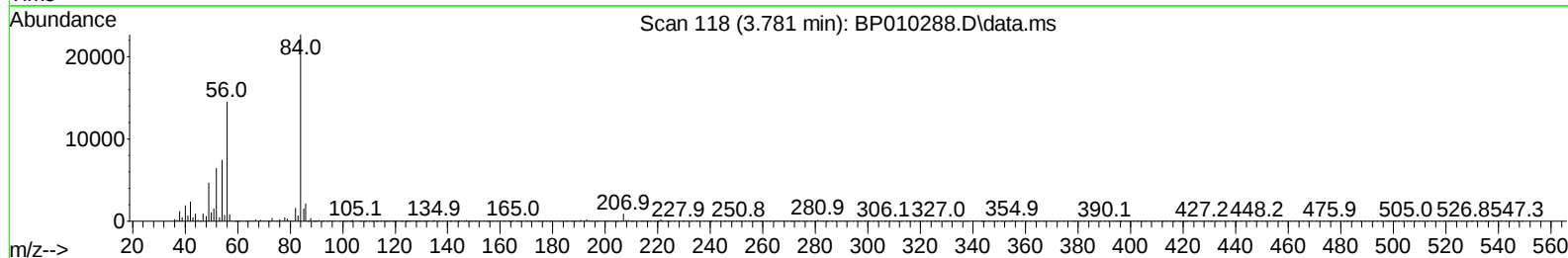
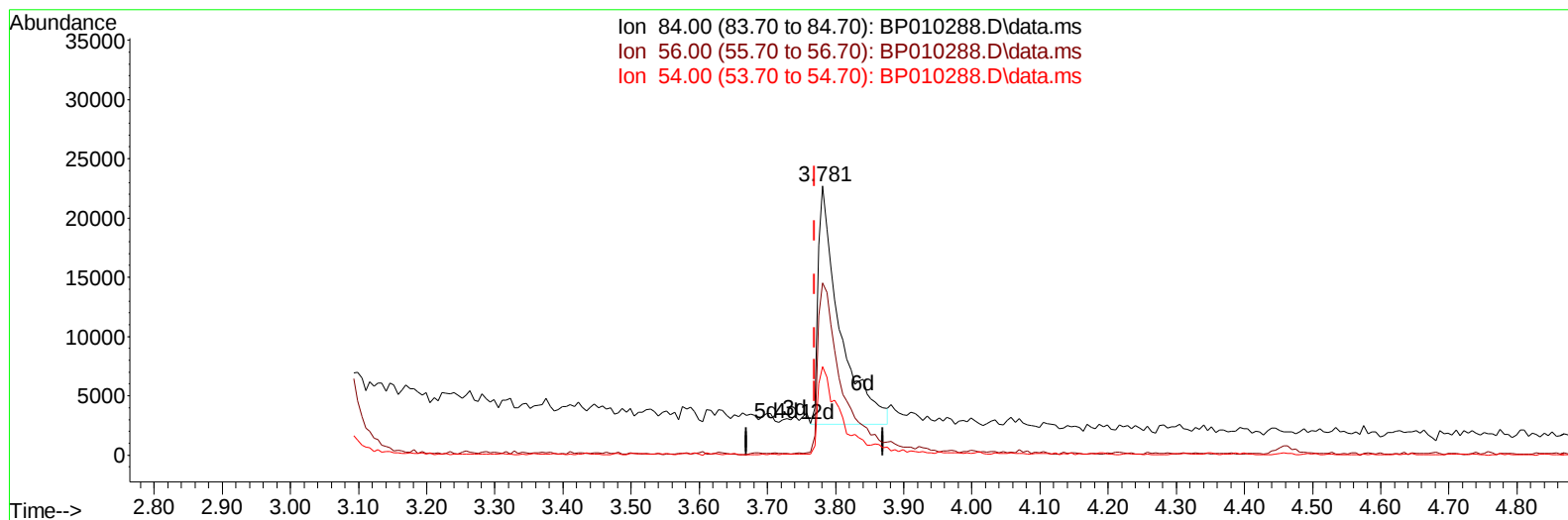
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(4) Pyridine-d5 (S)

3.781min (+ 0.012) 5.27 ng/ul m

response 43512

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	30.50	64.00#
54.00	24.80	32.80#
0.00	0.00	0.00

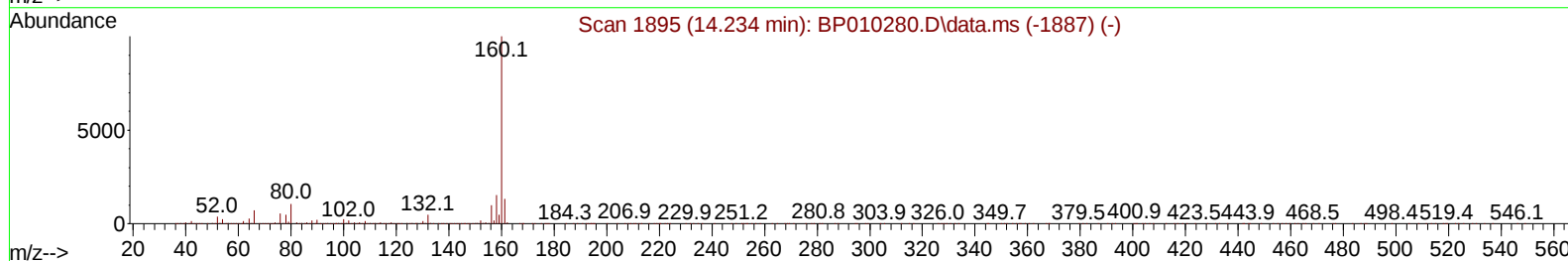
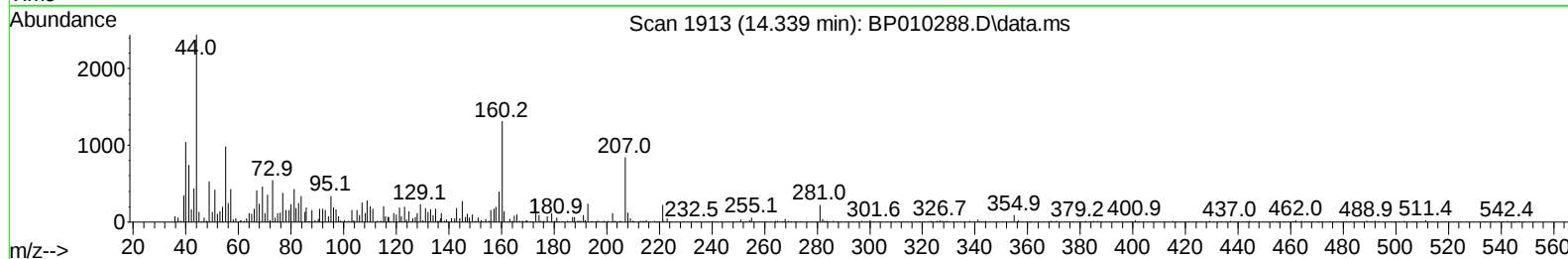
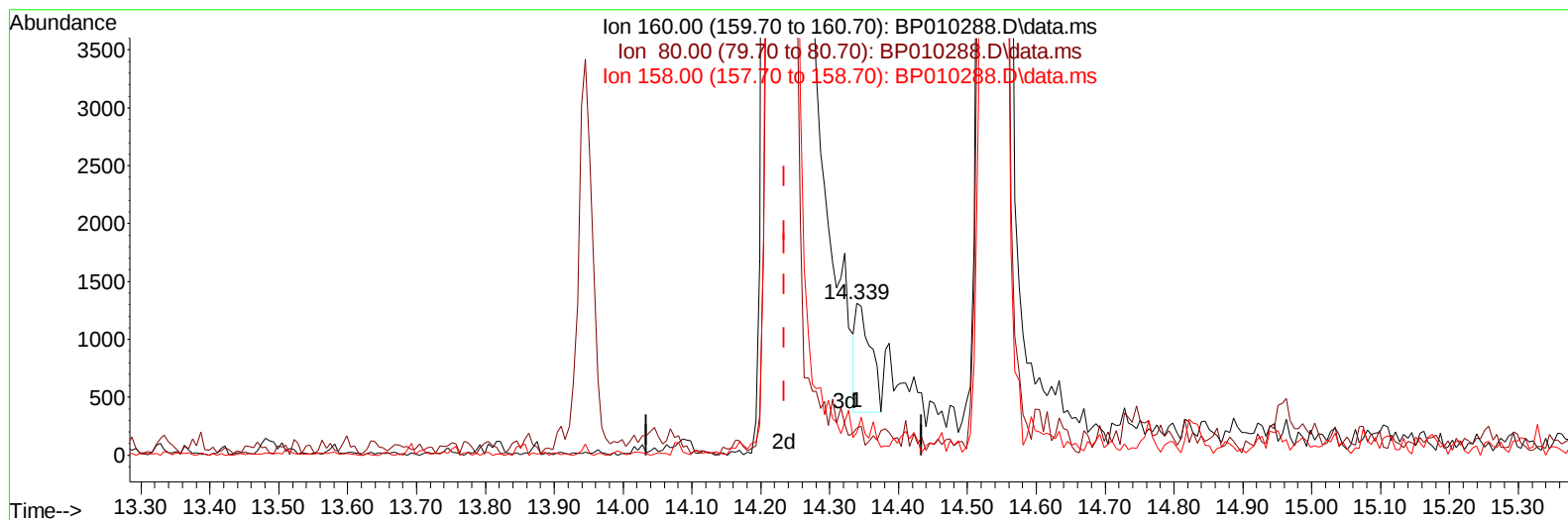
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Instrument :
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TIC: BP010288.D\data.ms

(49) Acenaphthylene-d8 (S)

14.339min (+ 0.106) 0.04 ng/u1

response 1420

Ion	Exp%	Act%
160.00	100.00	100.00
80.00	17.00	17.94
158.00	13.10	15.11
0.00	0.00	0.00

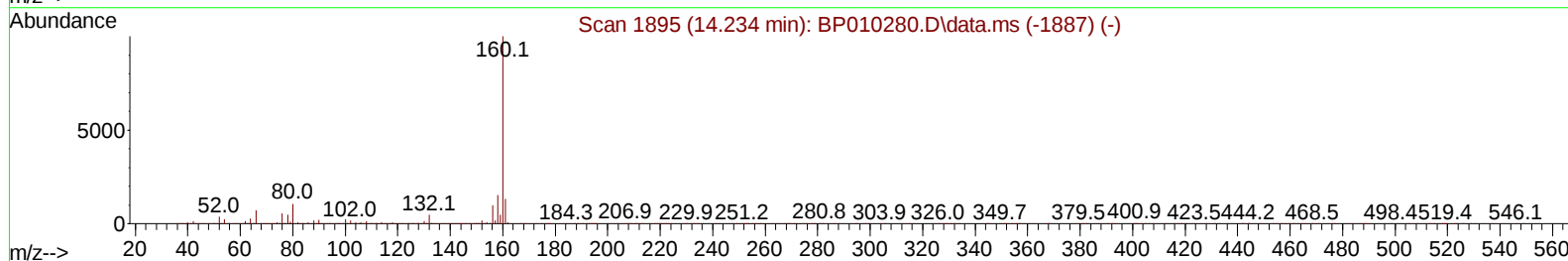
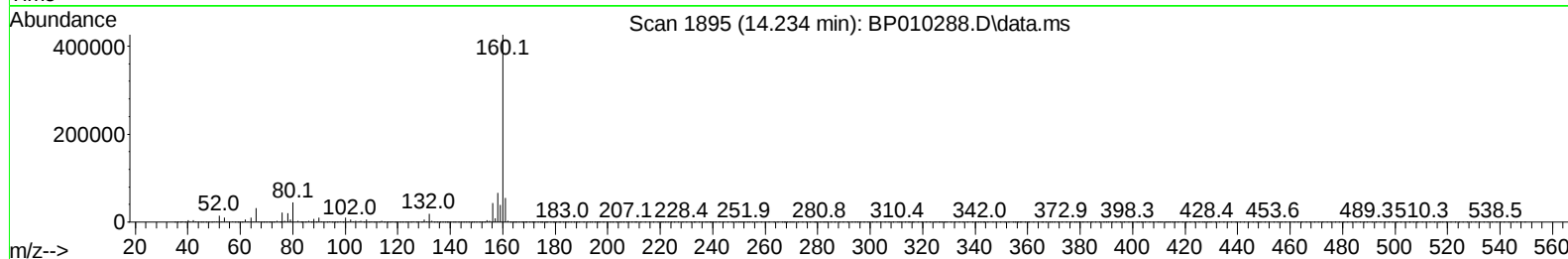
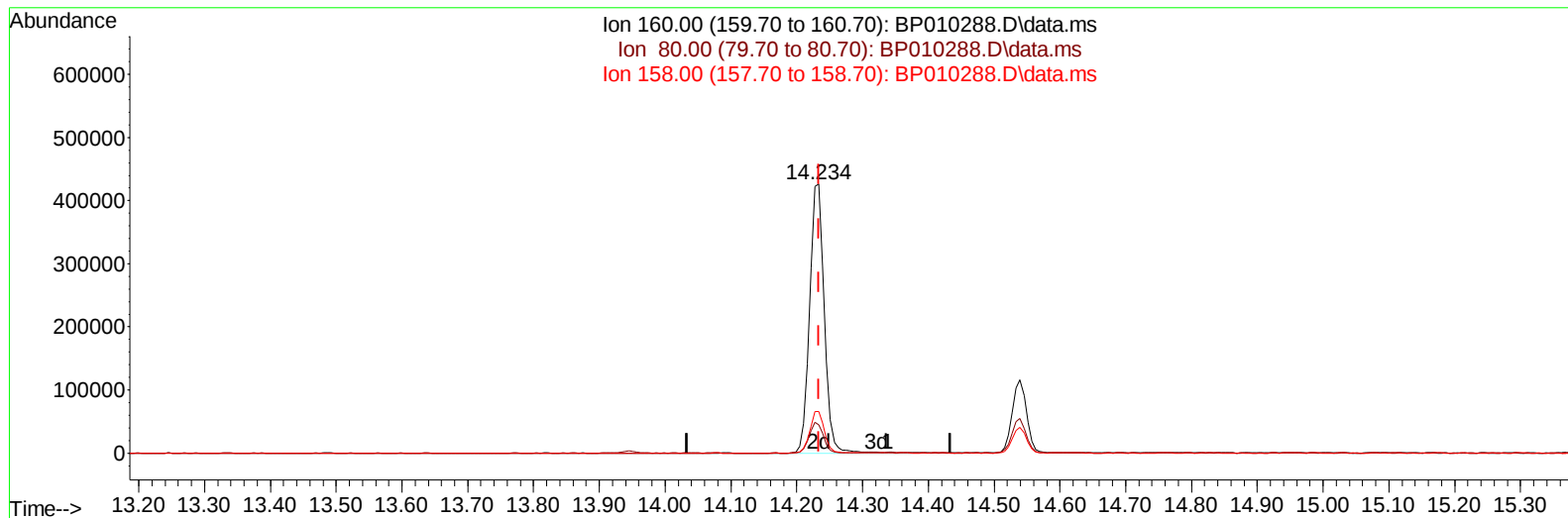
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Instrument :
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TIC: BP010288.D\data.ms

(49) Acenaphthylene-d8 (S)

14.234min (0.000) 17.38 ng/ul m

response 669468

Ion	Exp%	Act%
160.00	100.00	100.00
80.00	17.00	10.51#
158.00	13.10	15.54
0.00	0.00	0.00

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Instrument :
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Manual IntegrationsAPPROVED

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.905	152	125995	20.000	ng/ul	0.00
20) Naphthalene-d8	10.710	136	563711	20.000	ng/ul	# 0.00
38) Acenaphthene-d10	14.539	164	405485	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.286	188	899618	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.374	240	905014	20.000	ng/ul	# 0.00
88) Perylene-d12	23.804	264	932416	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.352	96	10060	3.392	ng/uL	0.00
4) Pyridine-d5	3.781	84	43512m	5.272	ng/ul	0.01
7) Phenol-d5	7.069	99	184657	15.780	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.234	67	134377	18.562	ng/ul	0.00
11) 2-Chlorophenol-d4	7.440	132	153964	17.865	ng/ul	0.00
15) 4-Methylphenol-d8	8.616	113	161884	16.826	ng/ul	0.00
21) Nitrobenzene-d5	9.063	128	81454	18.021	ng/ul	0.00
24) 2-Nitrophenol-d4	9.793	143	85128	16.519	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.334	165	150504	16.045	ng/ul	0.00
31) 4-Chloroaniline-d4	10.846	131	209454	15.126	ng/ul	0.00
46) Dimethylphthalate-d6	13.945	166	610738	18.906	ng/ul	0.00
49) Acenaphthylene-d8	14.234	160	669468m	17.382	ng/ul	0.00
54) 4-Nitrophenol-d4	14.739	143	60578	9.776	ng/ul	0.00
60) Fluorene-d10	15.534	176	527589	19.348	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.651	200	61229	9.939	ng/ul	0.00
73) Anthracene-d10	17.386	188	856283	19.437	ng/ul	0.00
81) Pyrene-d10	19.622	212	1114356	22.282	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.651	264	1066751	21.453	ng/ul	0.00
Target Compounds						
					Qvalue	
16) Acetophenone	8.734	105	15319	1.019	ng/ul#	74

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

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