

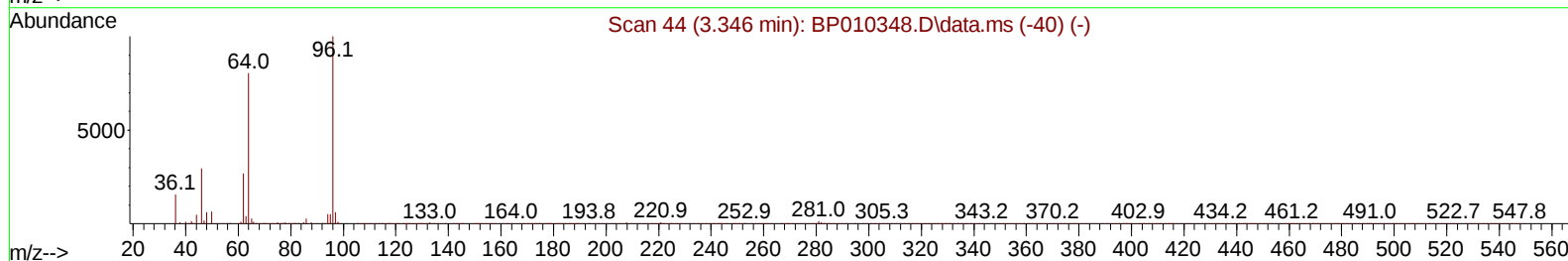
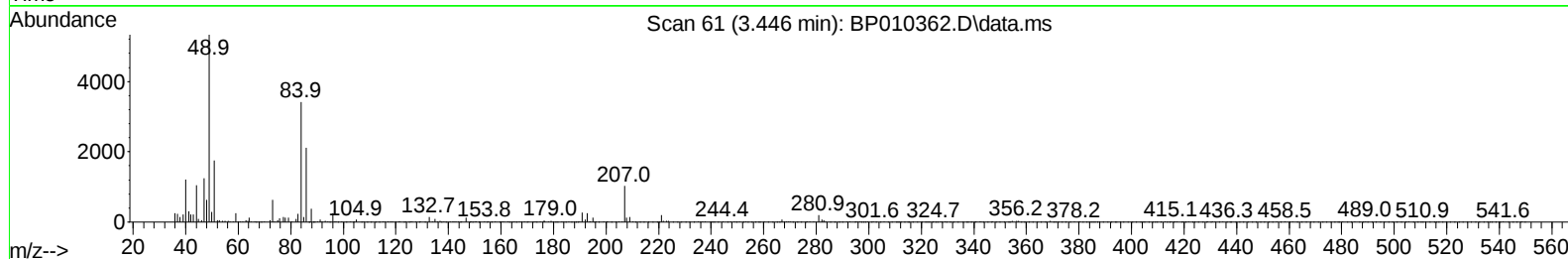
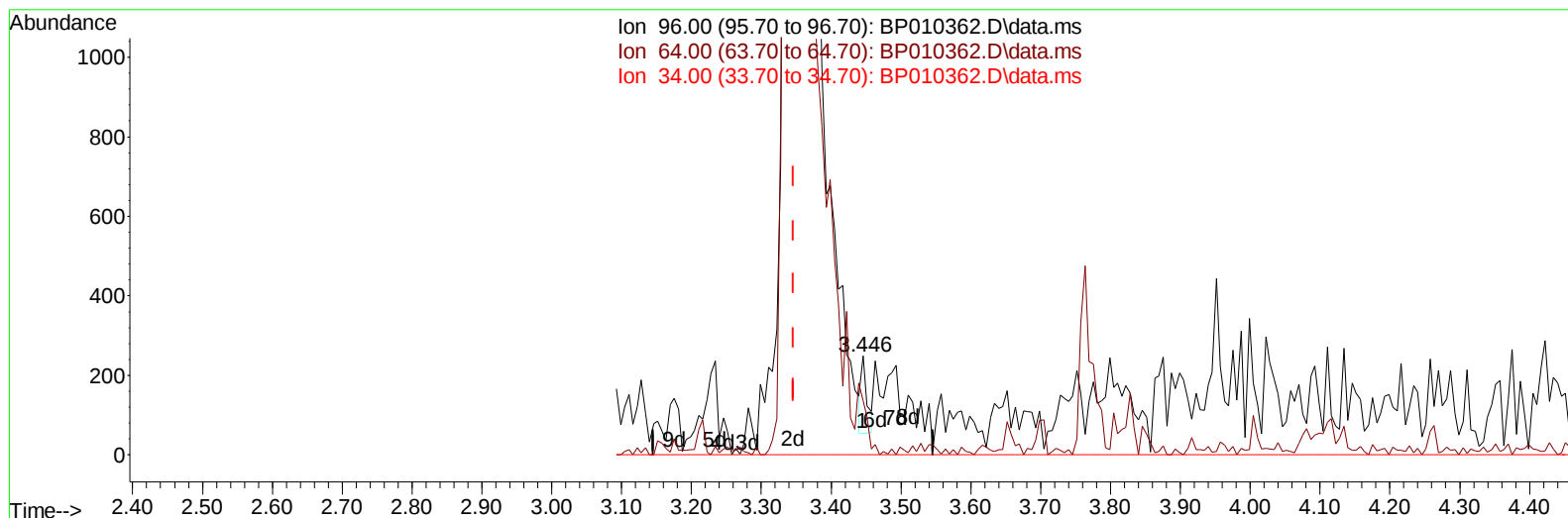
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051622\
 Data File : BP010362.D
 Acq On : 16 May 2022 18:04
 Operator : CG/JU
 Sample : N2713-15
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 DBSL4

Manual IntegrationsAPPROVED

Quant Time: May 16 23:21:48 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP051322.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri May 13 15:04:31 2022
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 05/17/2022
 Supervised By :mohammad ahmed 05/18/2022



TIC: BP010362.D\data.ms

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

3.446min (+ 0.100) 0.04 ng/uL

response 111

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	54.62
34.00	0.00	0.00
0.00	0.00	0.00

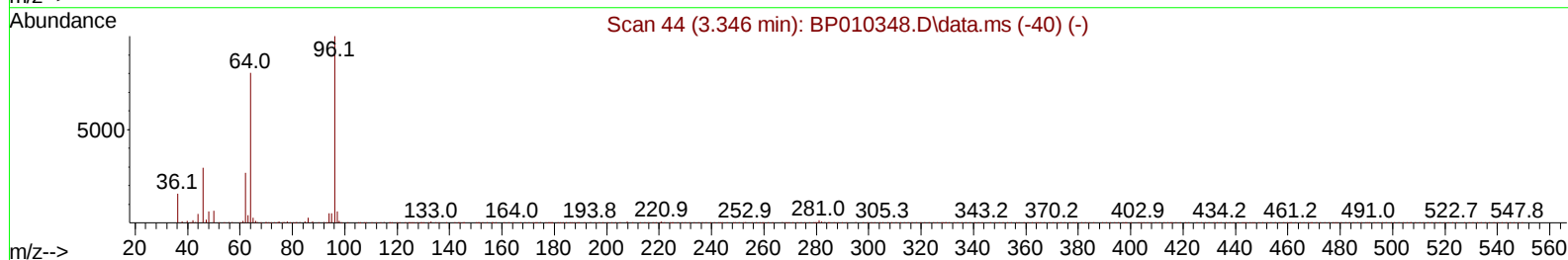
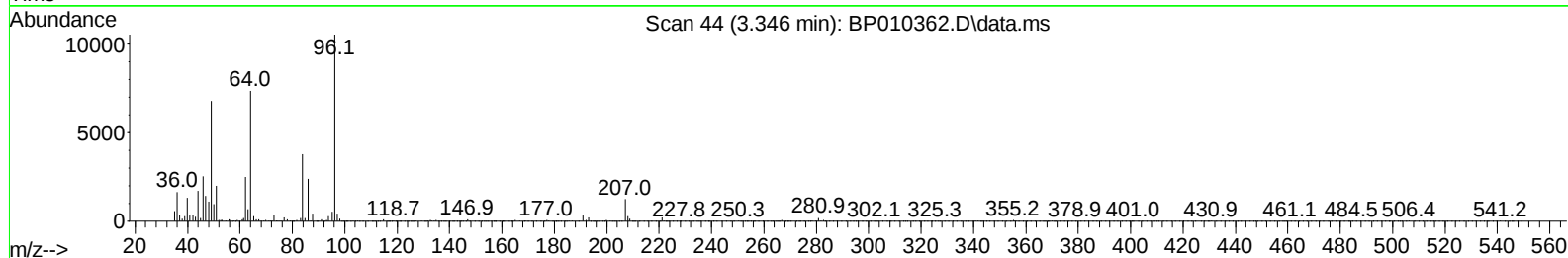
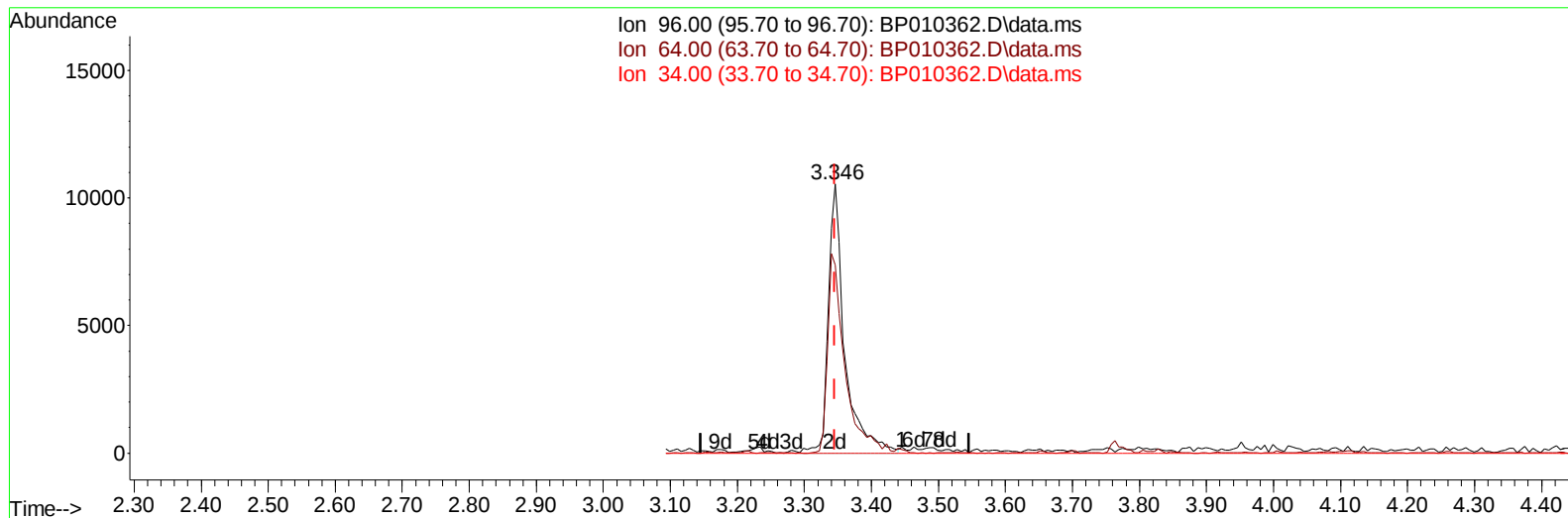
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TIC: BP010362.D\data.ms

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

3.346min (-0.000) 6.55 ng/uL m

response 17845

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	69.92#
34.00	0.00	0.00
0.00	0.00	0.00

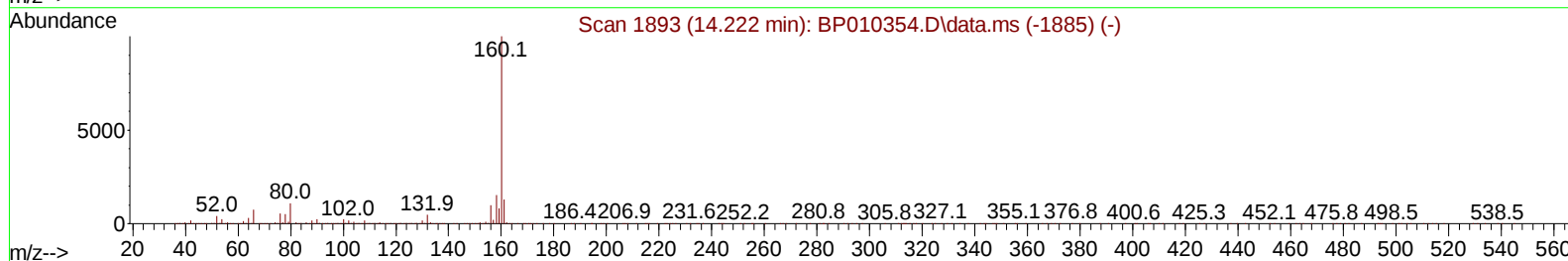
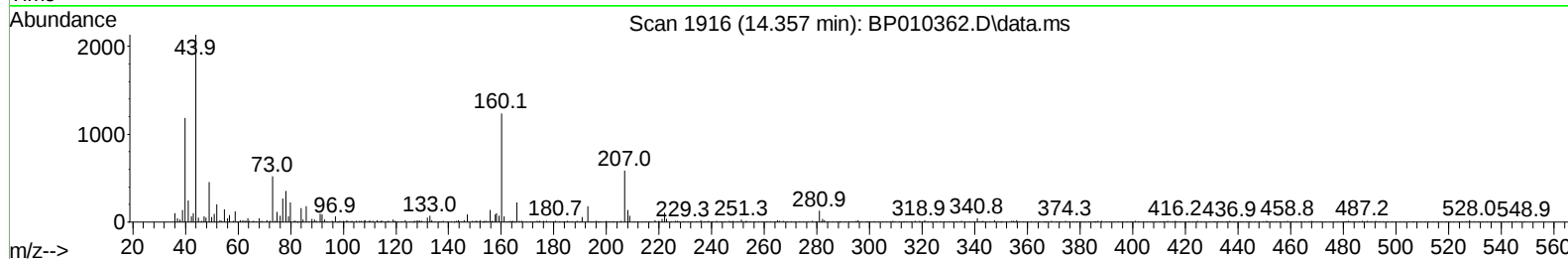
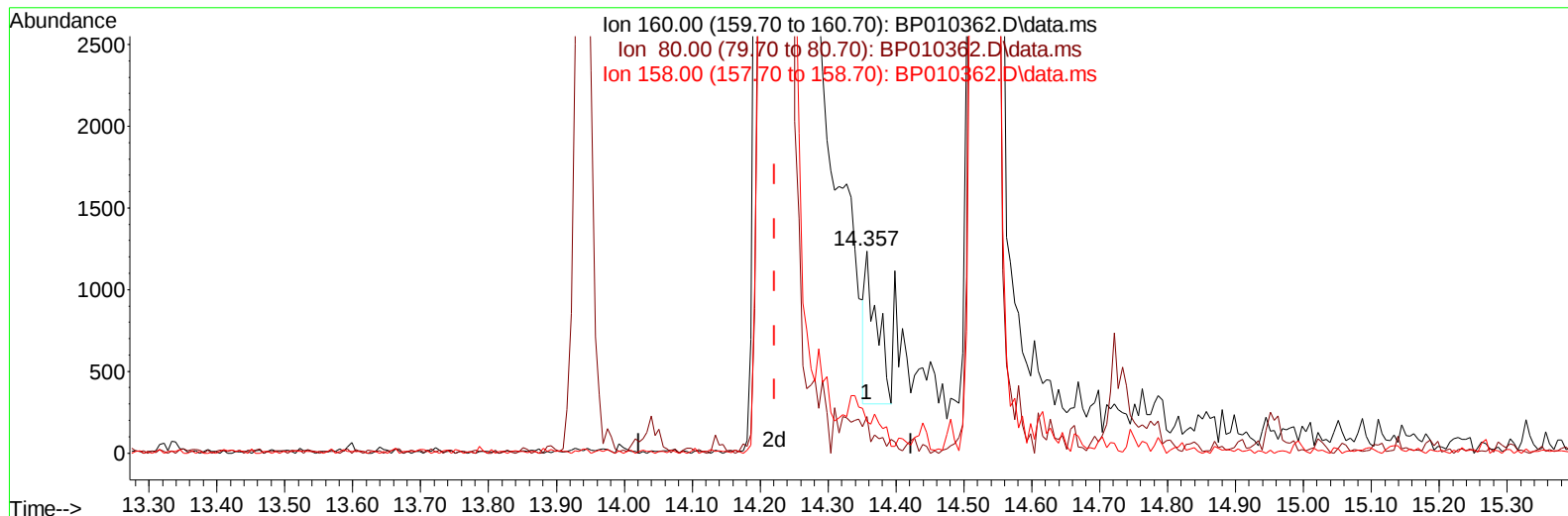
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TIC: BP010362.D\data.ms

(49) Acenaphthylene-d8 (S)

14.357min (+ 0.135) 0.04 ng/u1

response 1098

Ion	Exp%	Act%
160.00	100.00	100.00
80.00	17.00	17.95
158.00	13.10	15.44
0.00	0.00	0.00

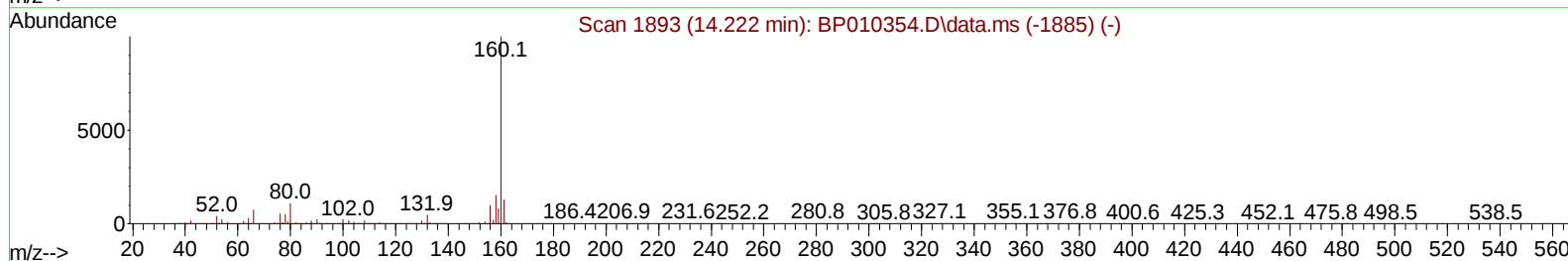
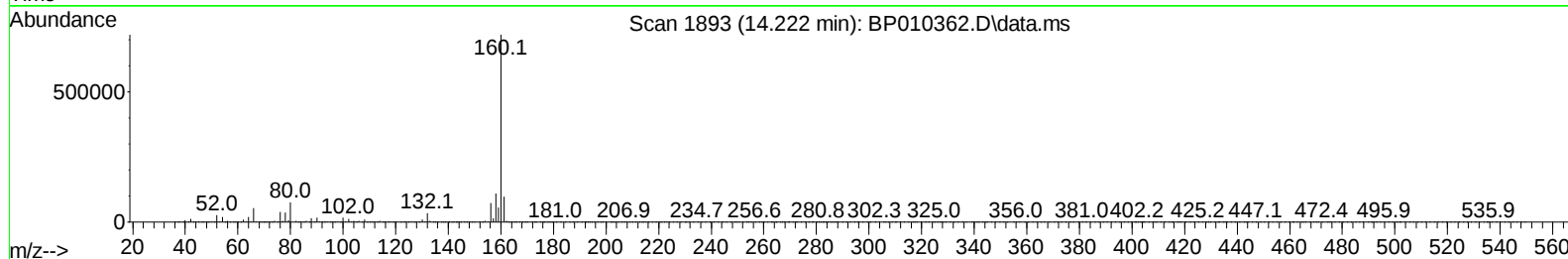
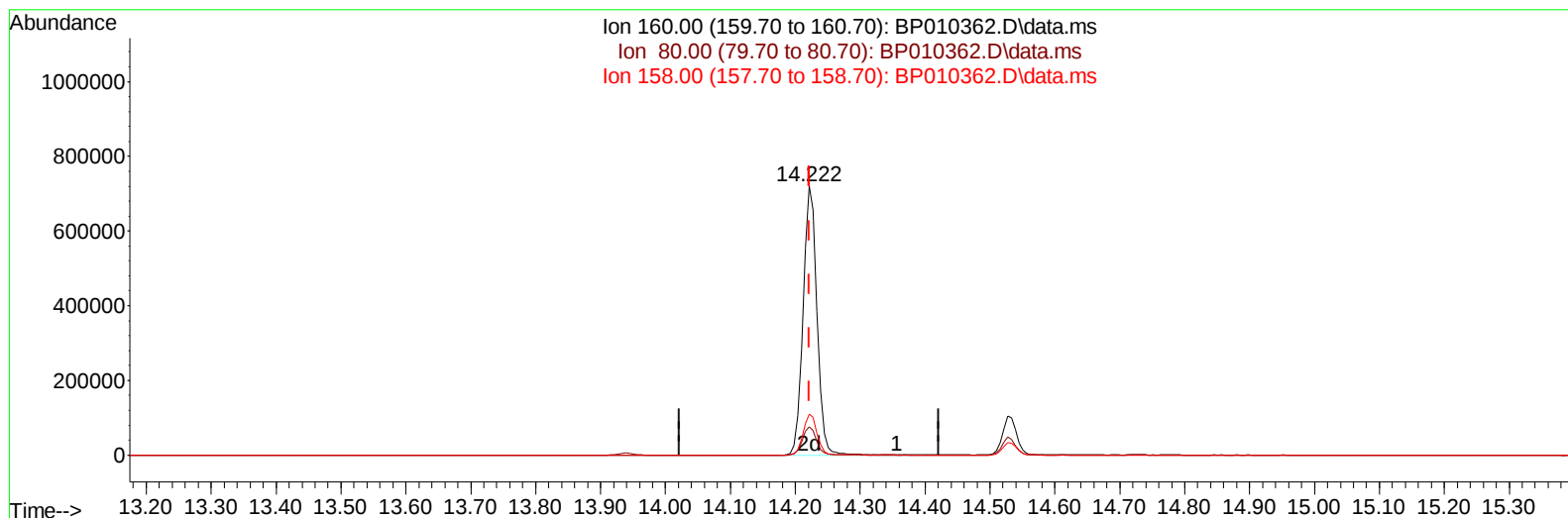
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TIC: BP010362.D\data.ms

(49) Acenaphthylene-d8 (S)

14.222min (-0.000) 35.63 ng/ul m

response 1072911

Ion	Exp%	Act%
160.00	100.00	100.00
80.00	17.00	10.57#
158.00	13.10	15.20
0.00	0.00	0.00

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 Misc :
 ALS Vial : 9 Sample Multiplier: 1

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	7.899	152	116115	20.000	ng/ul	0.00
20) Naphthalene-d8	10.698	136	504640	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.528	164	354380	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.280	188	791866	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.363	240	854433	20.000	ng/ul	# 0.00
88) Perylene-d12	23.792	264	866226	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.346	96	17845m	6.551	ng/uL	0.00
4) Pyridine-d5	3.763	84	232691	30.083	ng/ul	0.00
7) Phenol-d5	7.063	99	309547	31.762	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.228	67	222288	33.948	ng/ul	0.00
11) 2-Chlorophenol-d4	7.428	132	250108	33.538	ng/ul	0.00
15) 4-Methylphenol-d8	8.604	113	275024	33.206	ng/ul	0.00
21) Nitrobenzene-d5	9.057	128	136418	34.929	ng/ul	0.00
24) 2-Nitrophenol-d4	9.781	143	143425	34.652	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.322	165	256698	32.800	ng/ul	0.00
31) 4-Chloroaniline-d4	10.834	131	383335	32.942	ng/ul	0.00
46) Dimethylphthalate-d6	13.939	166	951680	36.449	ng/ul	0.00
49) Acenaphthylene-d8	14.222	160	1072911m	35.625	ng/ul	0.00
54) 4-Nitrophenol-d4	14.728	143	120146	25.853	ng/ul	0.00
60) Fluorene-d10	15.522	176	828254	36.463	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.645	200	121479	26.231	ng/ul	0.00
73) Anthracene-d10	17.380	188	1329490	37.078	ng/ul	0.00
81) Pyrene-d10	19.616	212	1664732	36.687	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.639	264	1611380	37.046	ng/ul	0.00

Target Compounds Qvalue

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

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