

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100322\
 Data File : BP011878.D
 Acq On : 03 Oct 2022 18:32
 Operator : CG/JU
 Sample : N4859-06
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

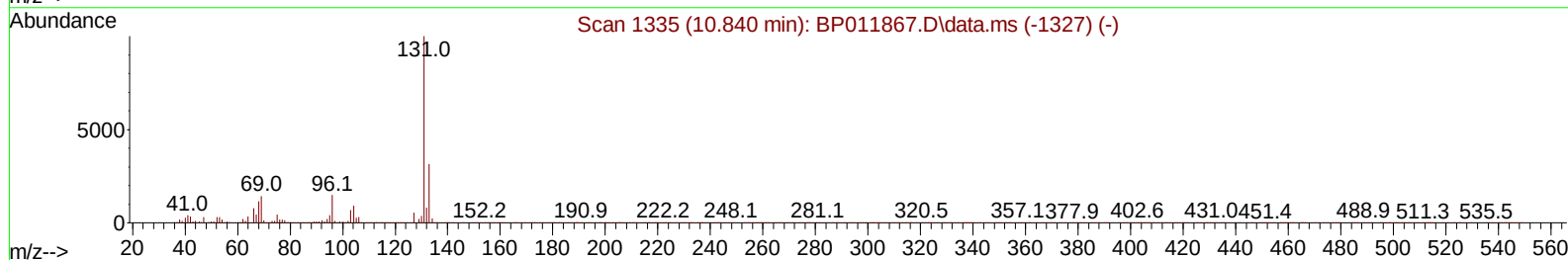
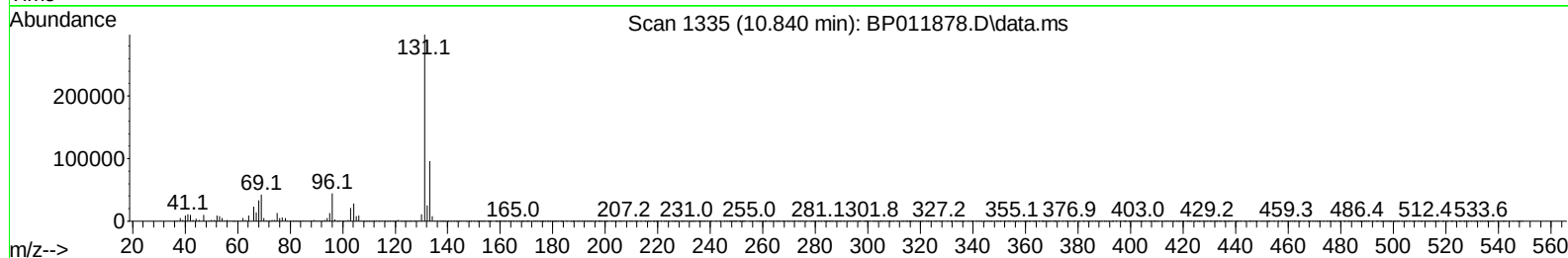
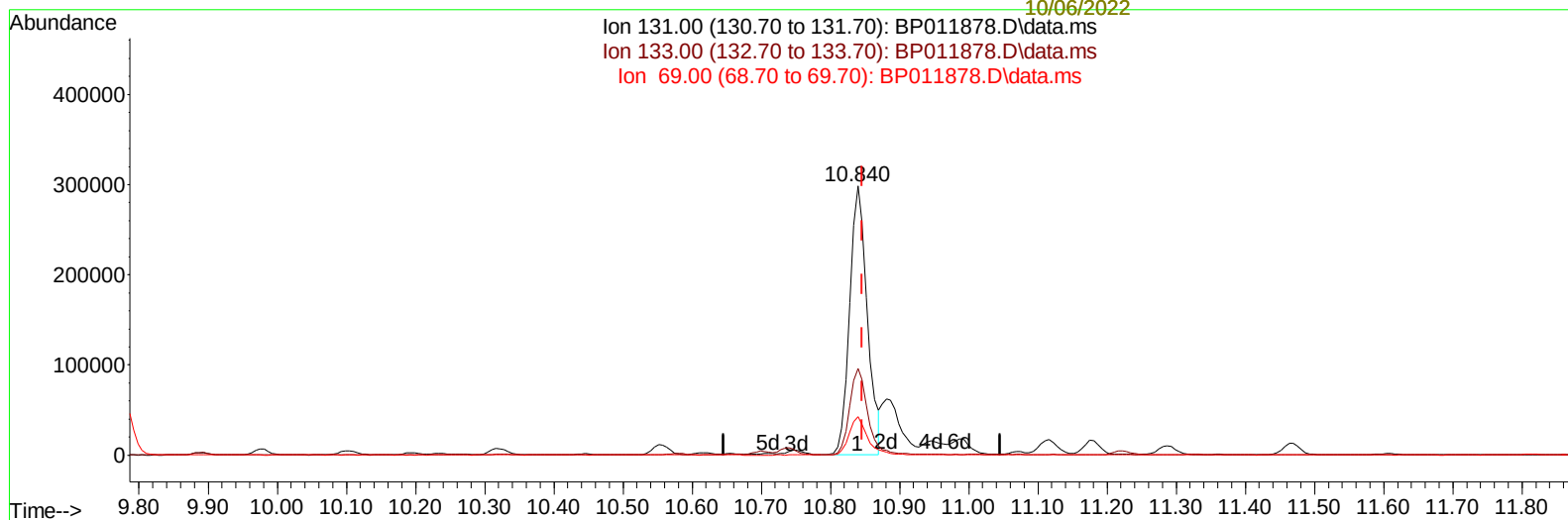
Instrument :
 BNA_P
 ClientSampleId :
 E10046

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 10/04/2022
 Supervised By :mohammad ahmed 10/06/2022

Supervised By :mohammad
 ahmed
 10/06/2022

Quant Time: Oct 03 23:32:50 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP093022.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Sat Oct 01 00:53:59 2022
 Response via : Initial Calibration



TIC: BP011878.D\data.ms

(31) 4-Chloroaniline-d4 (S)

10.840min (-0.006) 25.26 ng/ul

response 527272

Ion	Exp%	Act%
131.00	100.00	100.00
133.00	32.60	32.28
69.00	14.20	14.27
0.00	0.00	0.00

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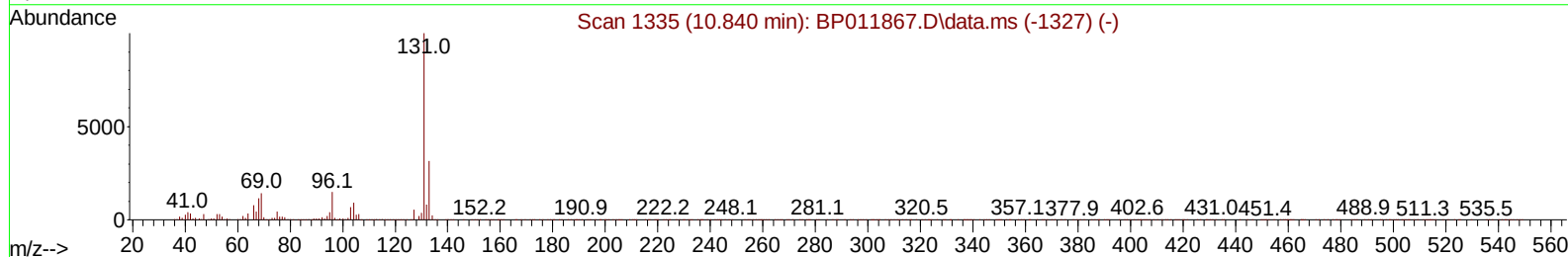
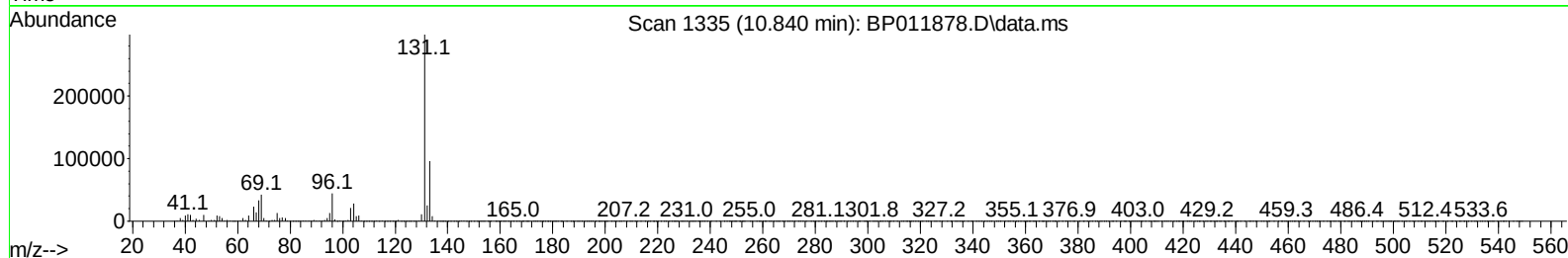
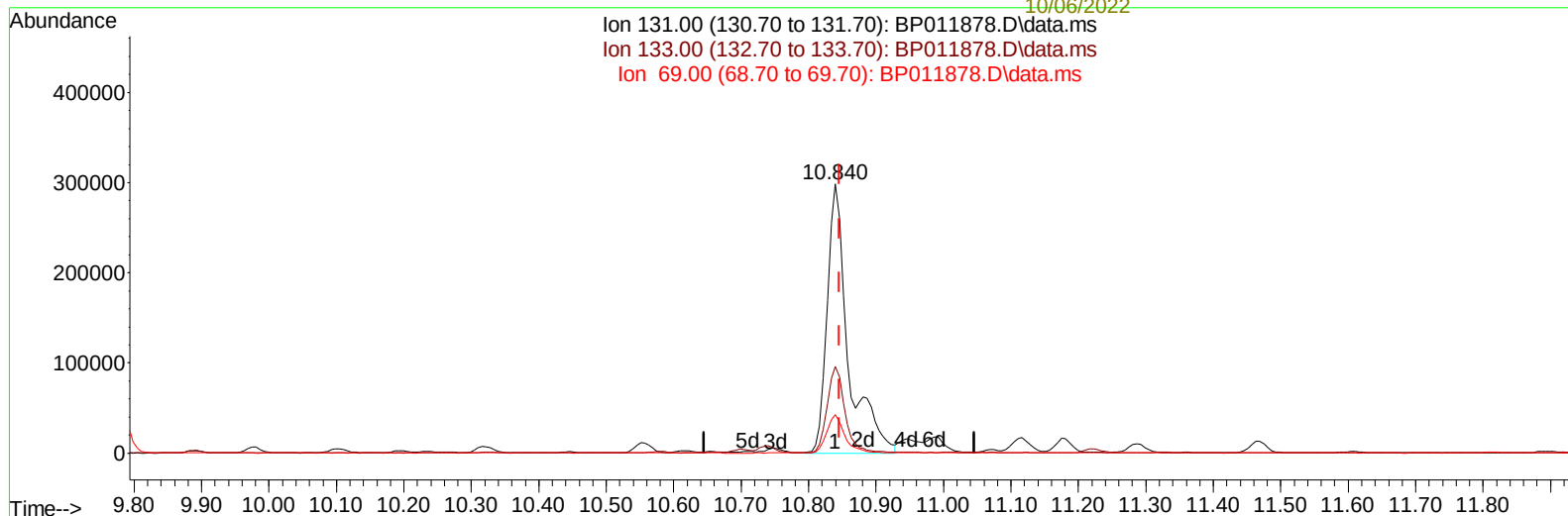
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(31) 4-Chloroaniline-d4 (S)

10.840min (-0.006) 31.03 ng/ul m

response 647851

Ion	Exp%	Act%
131.00	100.00	100.00
133.00	32.60	32.28
69.00	14.20	14.27
0.00	0.00	0.00

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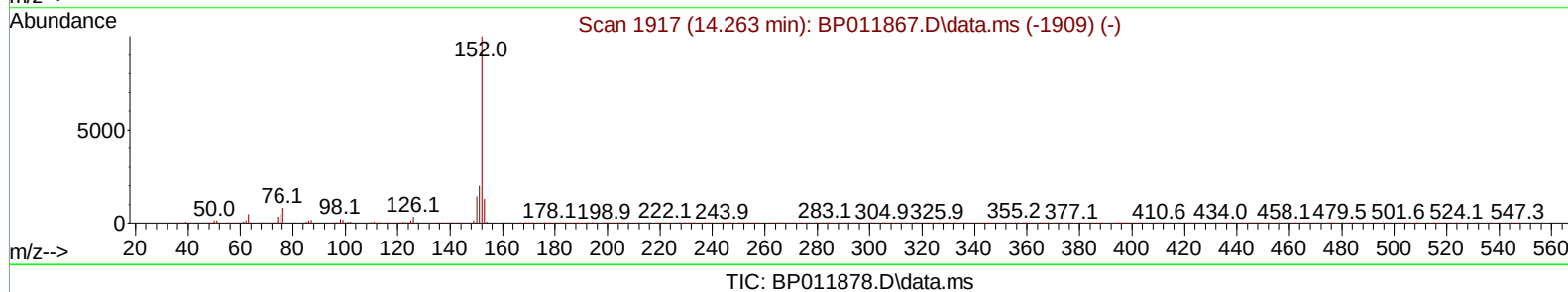
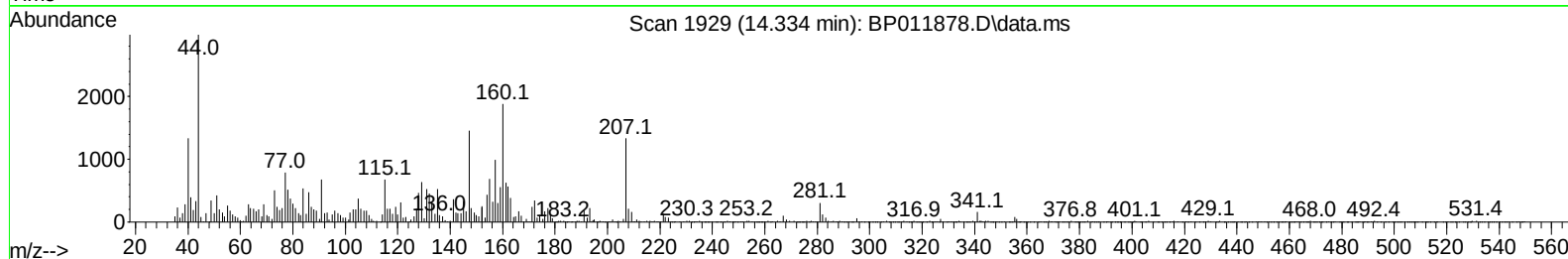
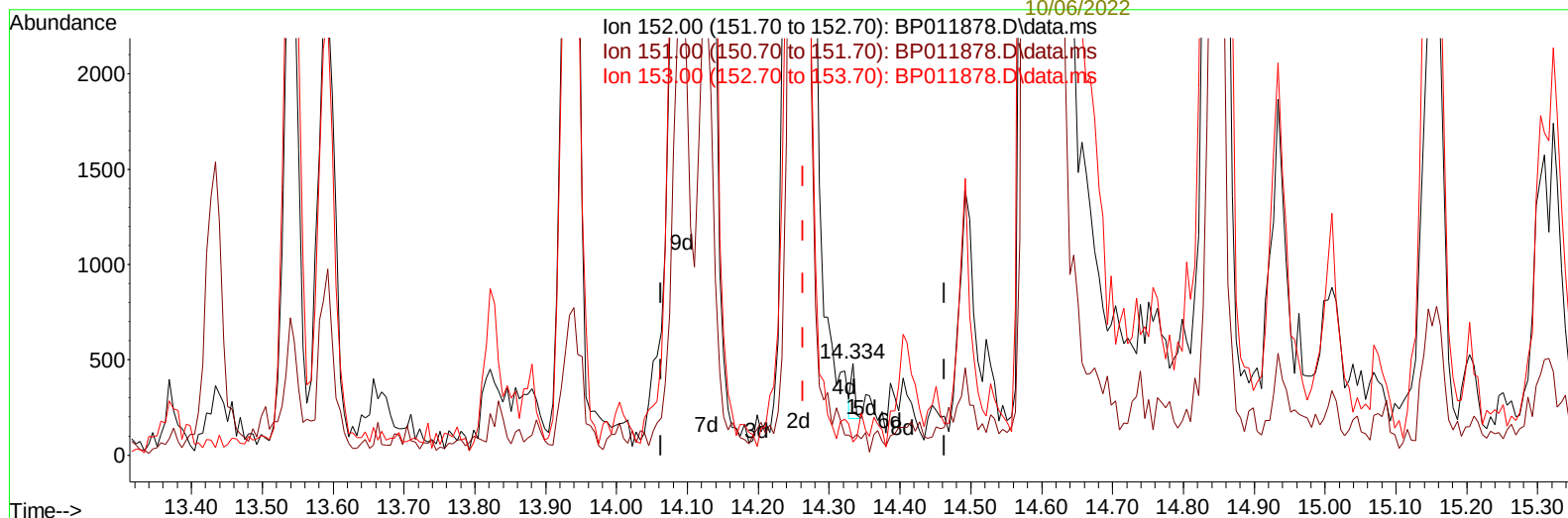
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(50) Acenaphthylene

14.334min (+ 0.070) 0.00 ng/u1

response	110	
Ion	Exp%	Act%
152.00	100.00	100.00
151.00	20.00	18.30
153.00	13.50	14.14
0.00	0.00	0.00

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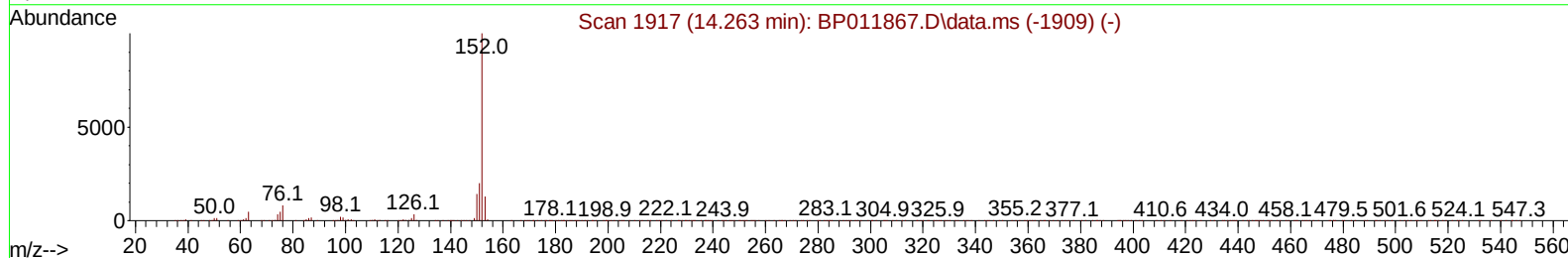
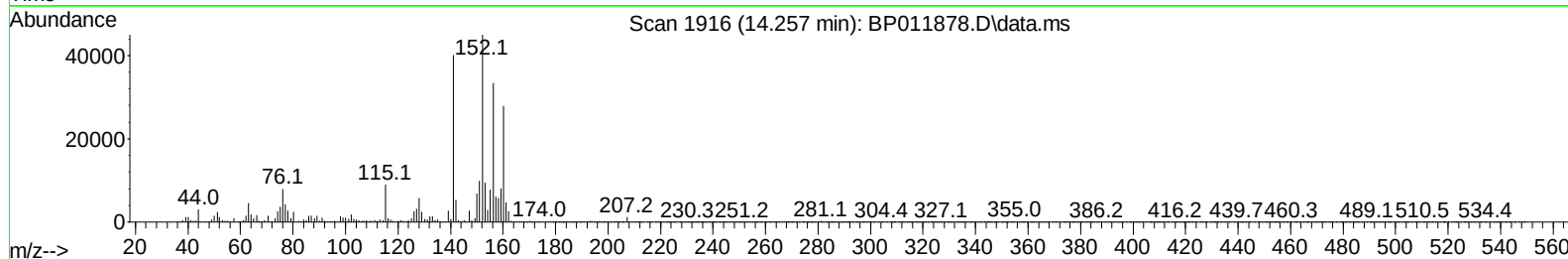
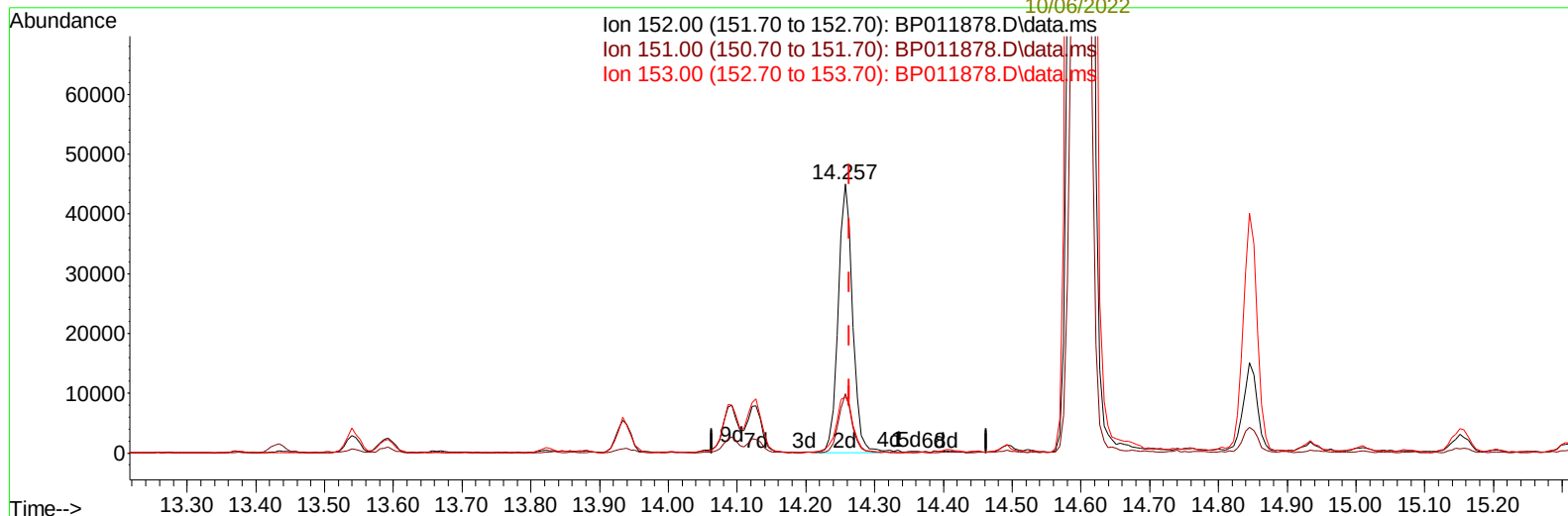
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(50) Acenaphthylene

14.257min (-0.006) 1.39 ng/ul m

response 65332

Ion	Exp%	Act%
152.00	100.00	100.00
151.00	20.00	22.12
153.00	13.50	20.97#
0.00	0.00	0.00

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.893	152	231711	20.000	ng/ul	0.00
20) Naphthalene-d8	10.698	136	908543	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.534	164	517968	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.292	188	1090393	20.000	ng/ul	0.00
79) Chrysene-d12	21.374	240	1200911	20.000	ng/ul	-0.01
88) Perylene-d12	23.810	264	1268921	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.305	96	26544	5.052	ng/uL	0.00
4) Pyridine-d5	3.728	84	134374	8.732	ng/ul	0.00
7) Phenol-d5	7.052	99	129284	6.463	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.222	67	328235	29.603	ng/ul	0.00
11) 2-Chlorophenol-d4	7.422	132	384755	24.366	ng/ul	0.00
15) 4-Methylphenol-d8	8.599	113	236274	14.288	ng/ul	0.00
21) Nitrobenzene-d5	9.057	128	229950	33.229	ng/ul	0.00
24) 2-Nitrophenol-d4	9.781	143	230767	31.168	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.316	165	397761	29.045	ng/ul	0.00
31) 4-Chloroaniline-d4	10.840	131	647851m	31.030	ng/ul	0.00
46) Dimethylphthalate-d6	13.945	166	1368823	36.955	ng/ul	0.00
49) Acenaphthylene-d8	14.228	160	1495169	35.409	ng/ul	0.00
54) 4-Nitrophenol-d4	14.739	143	47058	6.135	ng/ul	0.00
60) Fluorene-d10	15.528	176	1210390	37.628	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.645	200	193276	29.285	ng/ul	0.00
73) Anthracene-d10	17.392	188	2074199	41.760	ng/ul	0.00
81) Pyrene-d10	19.622	212	2576262	42.359	ng/ul	-0.01
92) Benzo(a)pyrene-d12	23.651	264	2777171	43.721	ng/ul	-0.01
Target Compounds						
30) Naphthalene	10.751	128	98994	2.029	ng/ul#	95
50) Acenaphthylene	14.257	152	65332m	1.393	ng/ul	
52) Acenaphthene	14.598	153	1897276	57.342	ng/ul	99
61) Fluorene	15.586	166	414233	11.160	ng/ul	100

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

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