

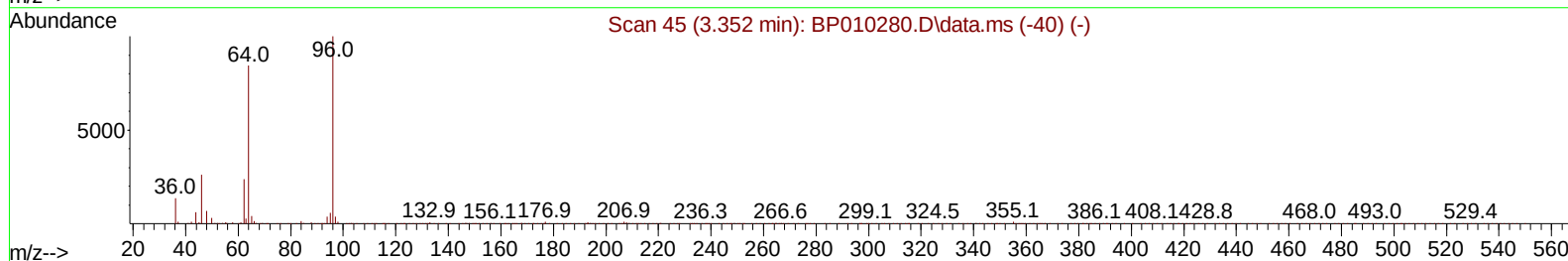
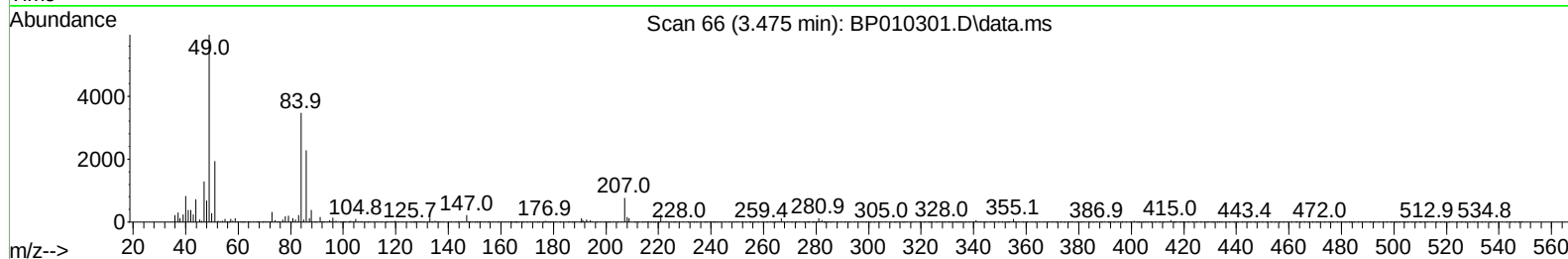
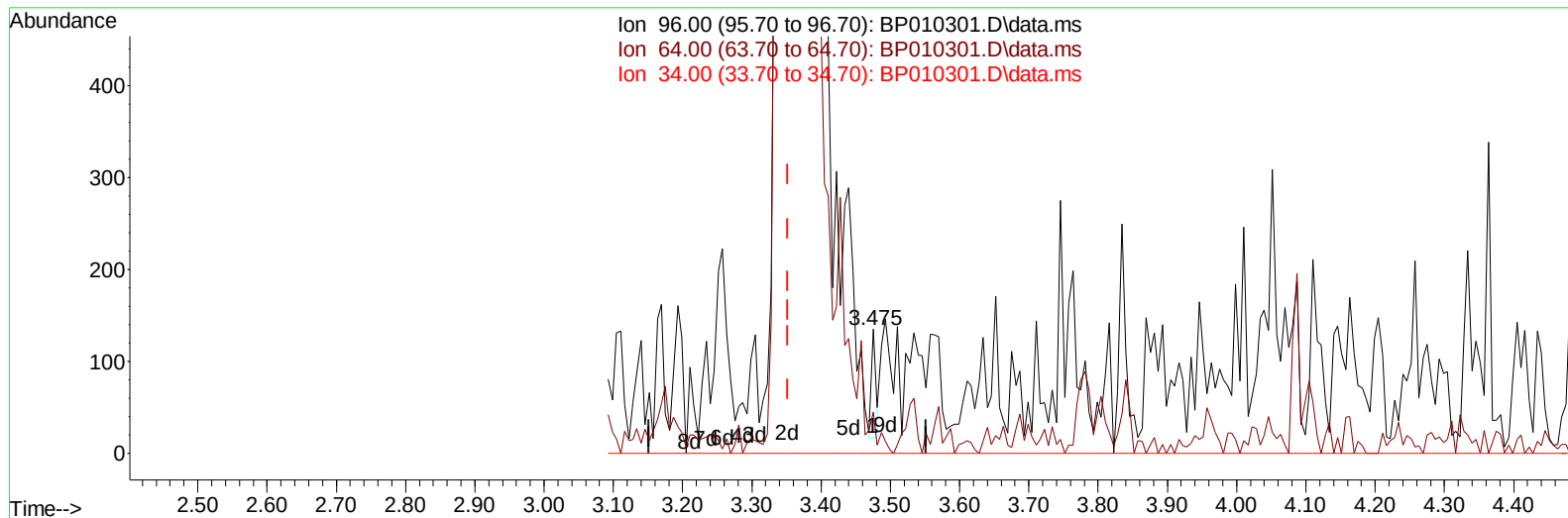
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051122\
 Data File : BP010301.D
 Acq On : 12 May 2022 00:28
 Operator : CG/JU
 Sample : N2707-09
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 EXL99

Manual IntegrationsAPPROVED

Quant Time: May 12 02:58:46 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: BP010301.D\data.ms

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

3.475min (+ 0.123) 0.02 ng/uL

response 54

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	45.93
34.00	0.00	0.00
0.00	0.00	0.00

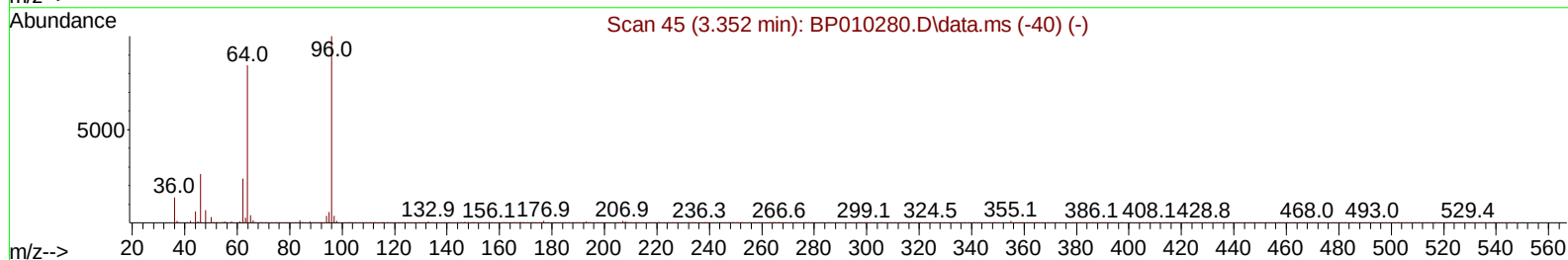
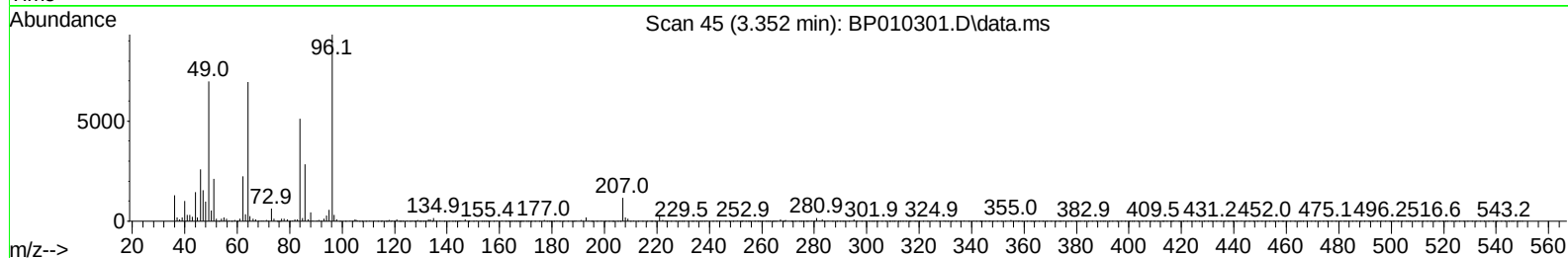
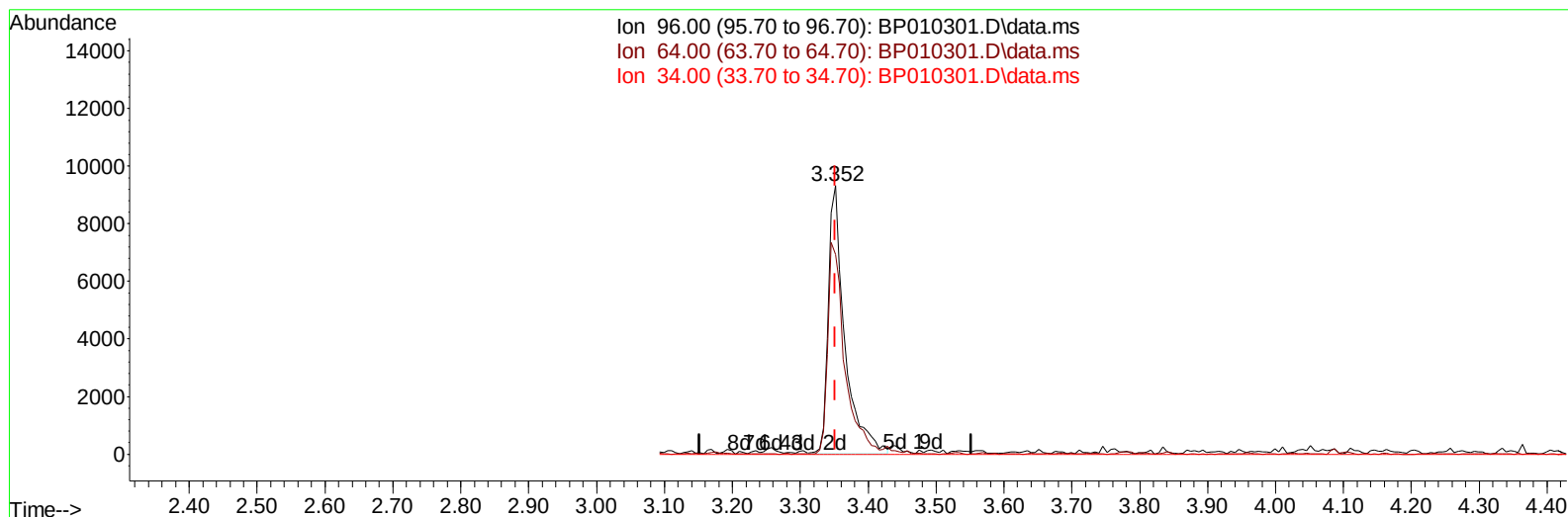
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TIC: BP010301.D\data.ms

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

3.352min (-0.000) 4.95 ng/uL m

response 15603

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	74.48#
34.00	0.00	0.00
0.00	0.00	0.00

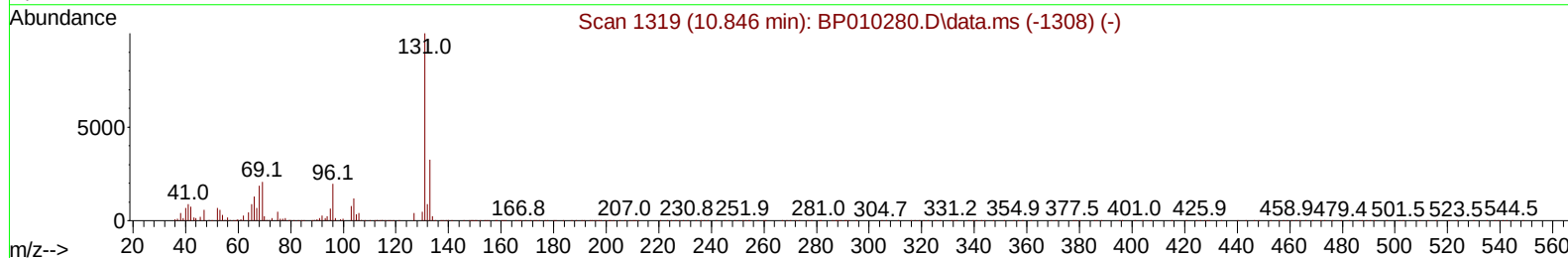
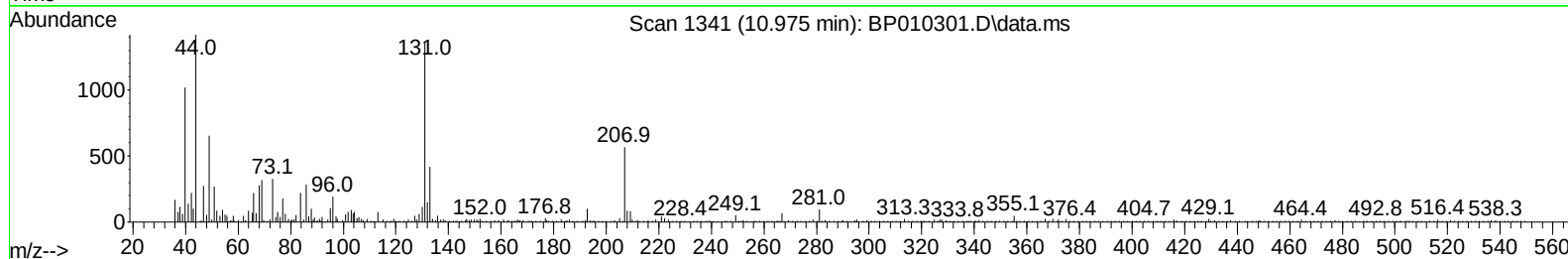
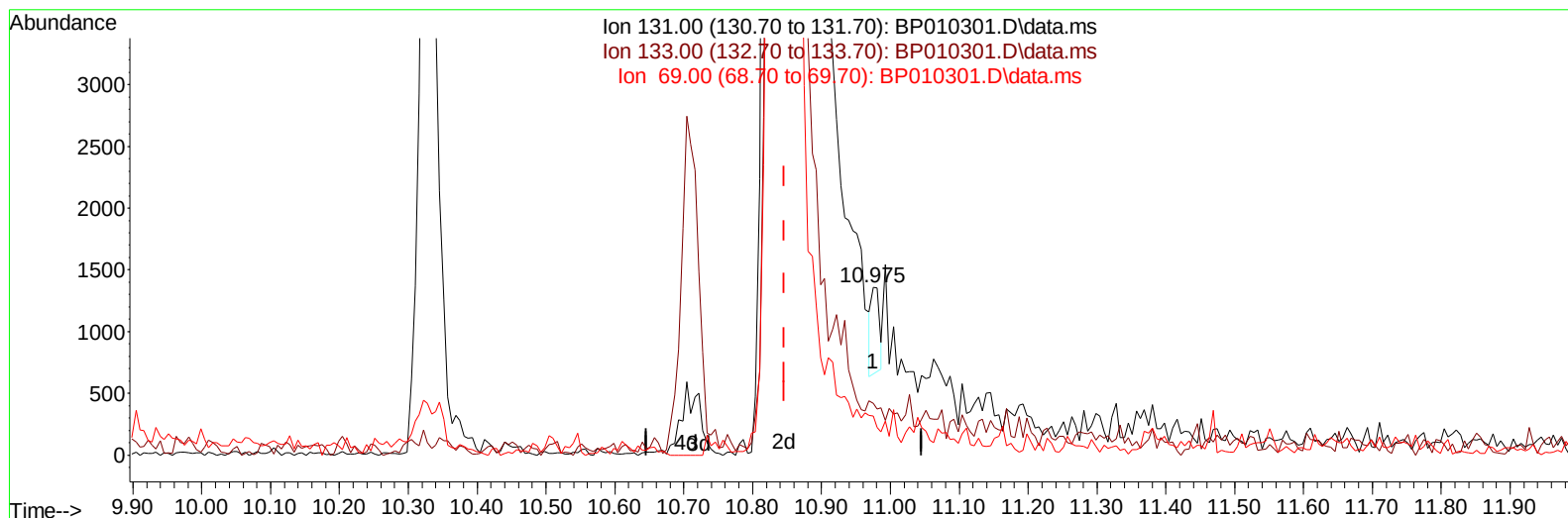
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Instrument :
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TIC: BP010301.D\data.ms

(31) 4-Chloroaniline-d4 (S)

10.975min (+ 0.129) 0.04 ng/u1

response 577

Ion	Exp%	Act%
131.00	100.00	100.00
133.00	29.20	30.74
69.00	26.10	23.38
0.00	0.00	0.00

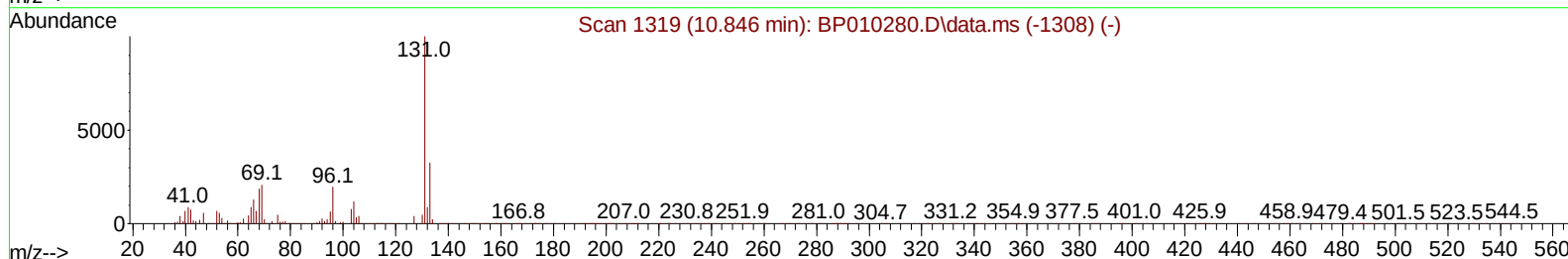
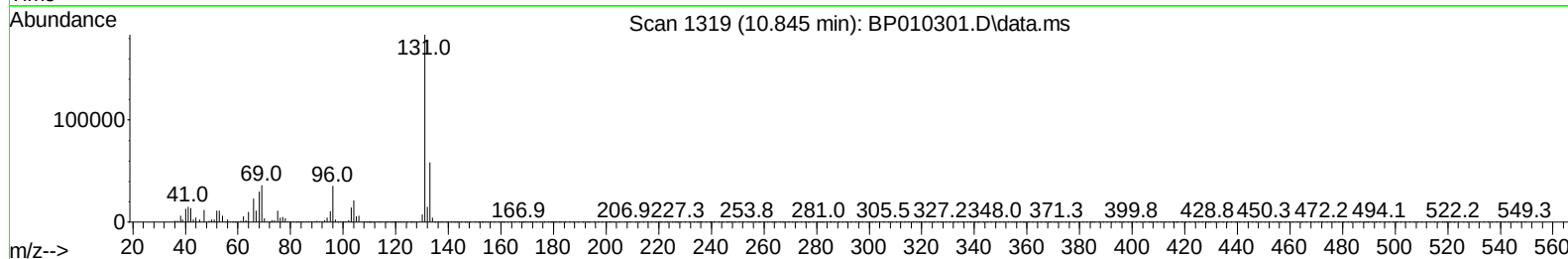
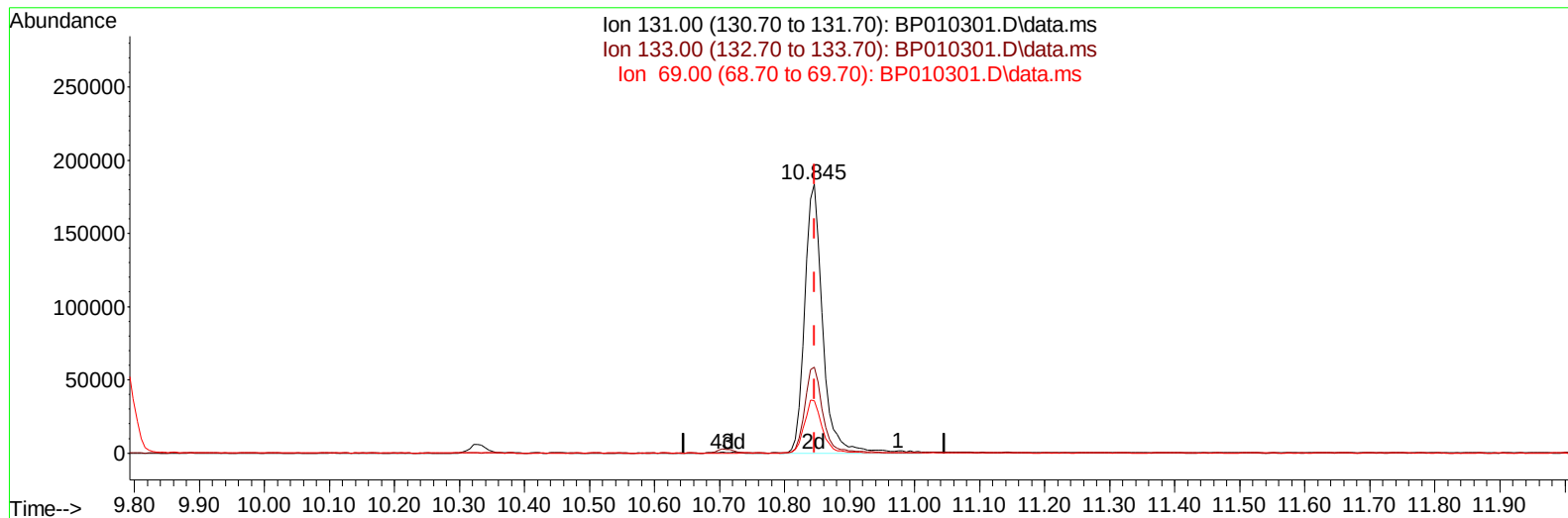
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Instrument :
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TIC: BP010301.D\data.ms

(31) 4-Chloroaniline-d4 (S)

10.845min (-0.000) 24.13 ng/ul m

response 355422

Ion	Exp%	Act%
131.00	100.00	100.00
133.00	29.20	31.98
69.00	26.10	19.58#
0.00	0.00	0.00

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.904	152	133947	20.000	ng/ul	0.00
20) Naphthalene-d8	10.704	136	599693	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.539	164	422096	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.286	188	934348	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.368	240	936073	20.000	ng/ul	# 0.00
88) Perylene-d12	23.798	264	933706	20.000	ng/ul	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.352	96	15603m	4.949	ng/uL	0.00
4) Pyridine-d5	3.781	84	39511	4.503	ng/ul	0.01
7) Phenol-d5	7.069	99	68862	5.535	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.234	67	267776	34.793	ng/ul	0.00
11) 2-Chlorophenol-d4	7.440	132	223381	24.380	ng/ul	0.00
15) 4-Methylphenol-d8	8.610	113	144457	14.124	ng/ul	0.00
21) Nitrobenzene-d5	9.063	128	163340	33.969	ng/ul	0.00
24) 2-Nitrophenol-d4	9.787	143	168132	30.667	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.328	165	272474	27.306	ng/ul	0.00
31) 4-Chloroaniline-d4	10.845	131	355422m	24.127	ng/ul	0.00
46) Dimethylphthalate-d6	13.945	166	1168495	34.749	ng/ul	0.00
49) Acenaphthylene-d8	14.228	160	1287550	32.115	ng/ul	0.00
54) 4-Nitrophenol-d4	14.739	143	26903	4.171	ng/ul	0.00
60) Fluorene-d10	15.528	176	1031914	36.354	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.651	200	197767	30.909	ng/ul	0.00
73) Anthracene-d10	17.386	188	1768991	38.662	ng/ul	0.00
81) Pyrene-d10	19.621	212	2175641	42.059	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.645	264	2059328	41.356	ng/ul	-0.01

Target Compounds Qvalue

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

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