

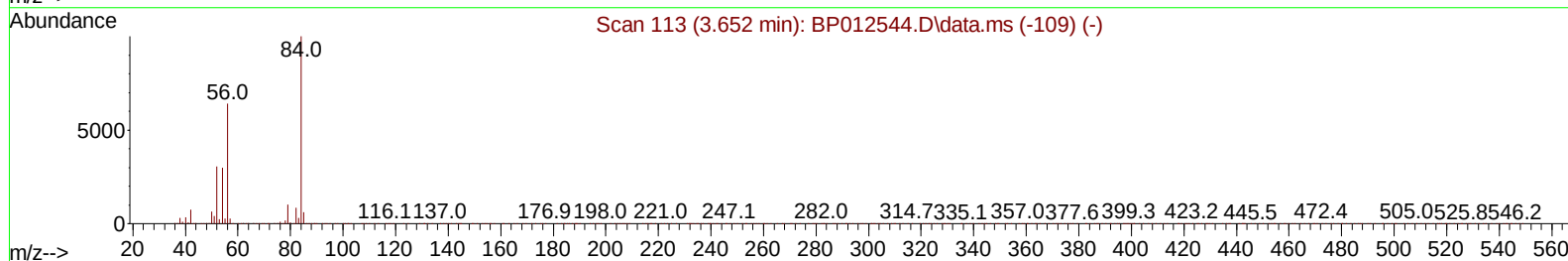
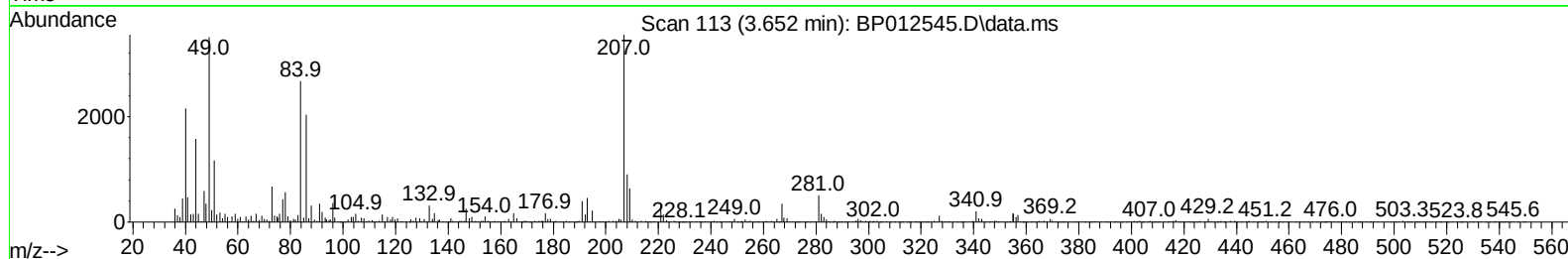
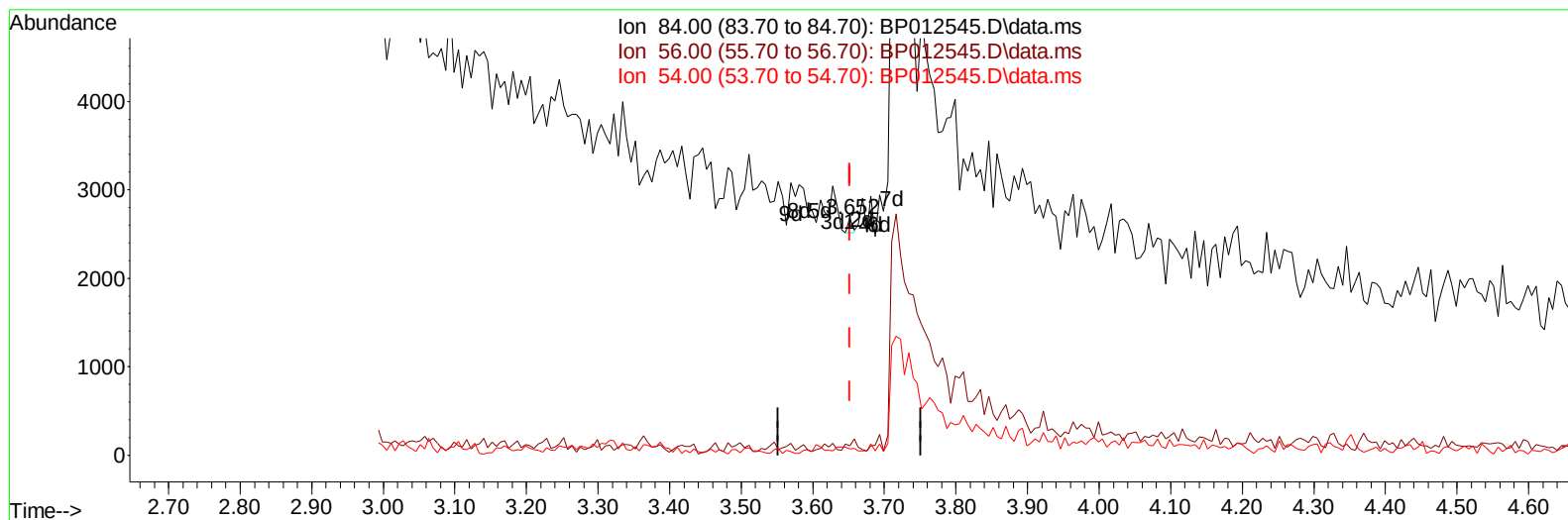
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110822\
 Data File : BP012545.D
 Acq On : 08 Nov 2022 10:38
 Operator : CG/JU
 Sample : PB148823BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK823

Manual IntegrationsAPPROVED

Quant Time: Nov 08 21:40:28 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Nov 04 02:53:36 2022
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 11/09/2022
 Supervised By :mohammad ahmed 11/14/2022



TIC: BP012545.D\data.ms

(4) Pyridine-d5 (S)

3.652min (-0.000) 0.00 ng/uI

response 67

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	65.20	4.30#
54.00	29.80	2.62#
0.00	0.00	0.00

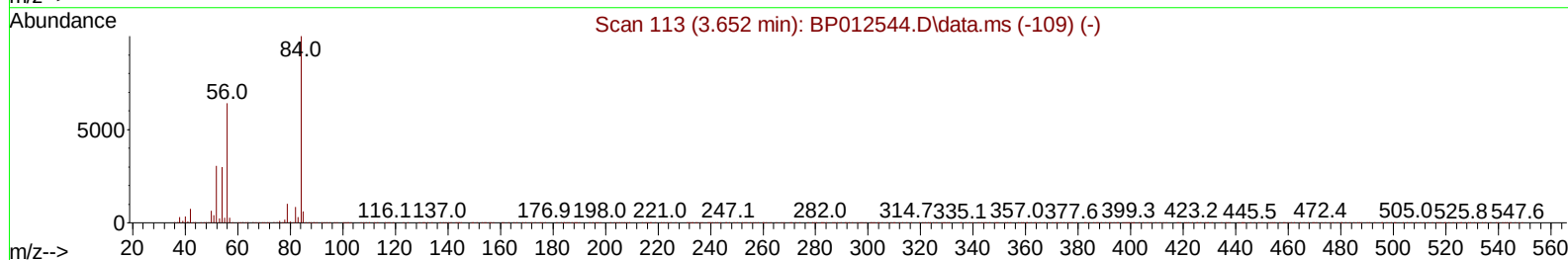
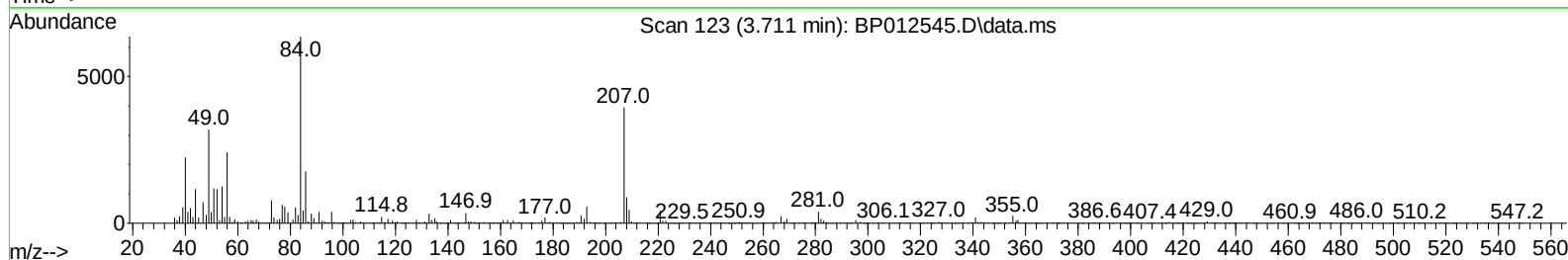
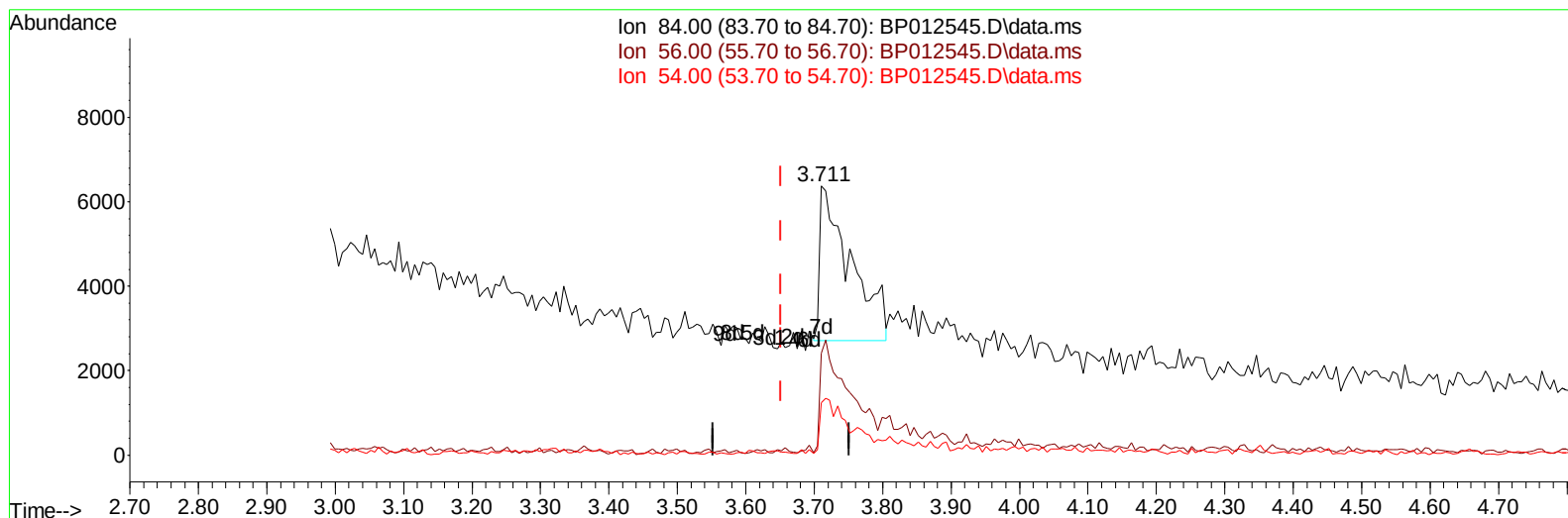
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(4) Pyridine-d5 (S)

3.711min (+ 0.059) 0.46 ng/ul m

response 11452

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	65.20	37.86#
54.00	29.80	19.58#
0.00	0.00	0.00

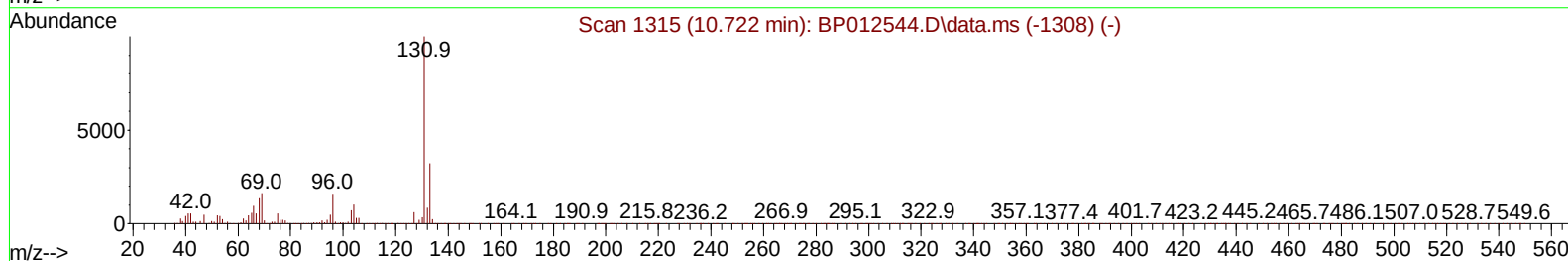
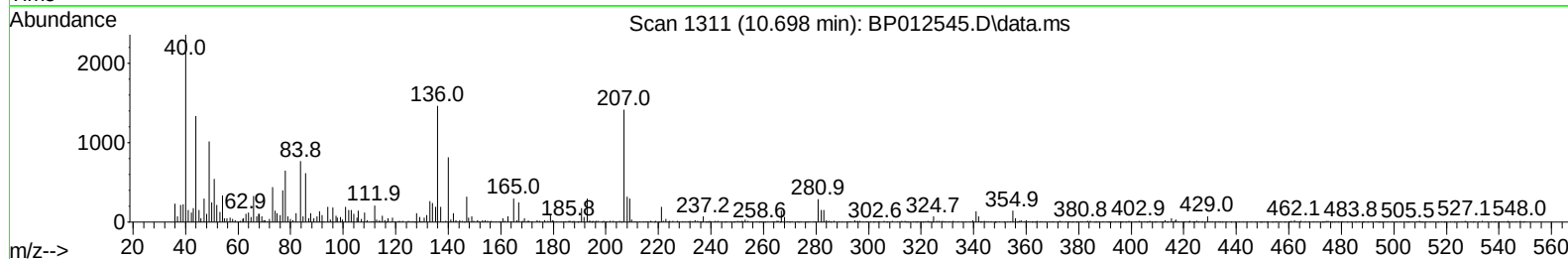
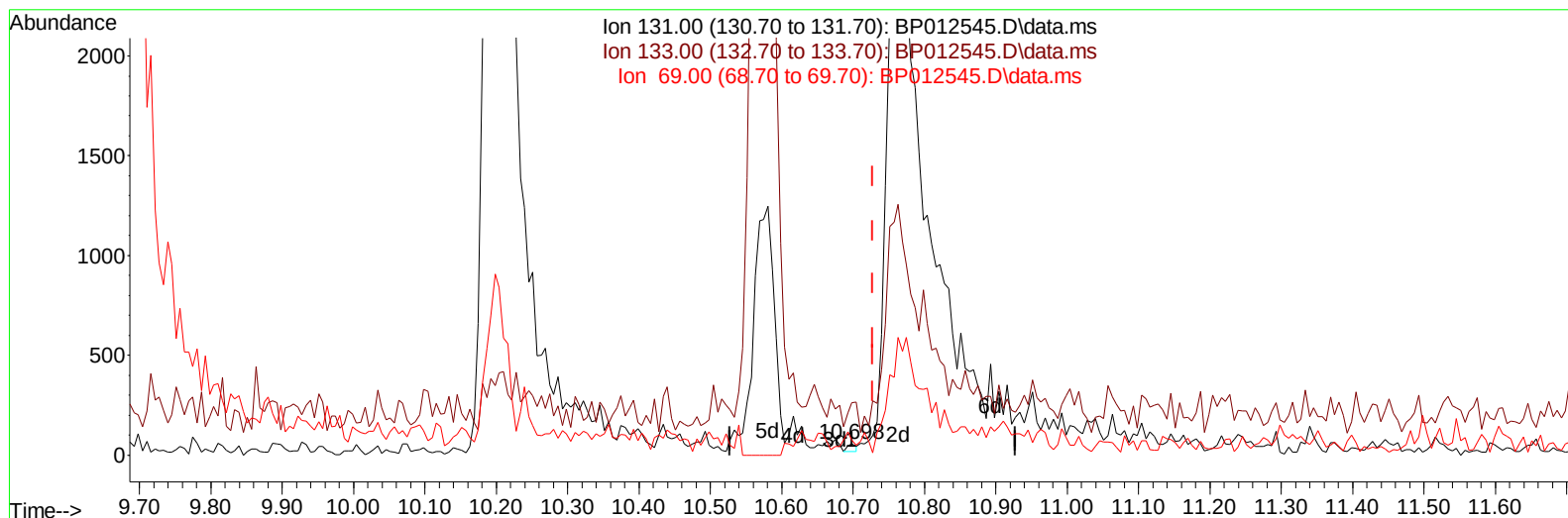
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TIC: BP012545.D\data.ms

(31) 4-Chloroaniline-d4 (S)

10.698min (-0.030) 0.00 ng/ul

response 32

Ion	Exp%	Act%
131.00	100.00	100.00
133.00	32.20	469.64#
69.00	16.50	133.93#
0.00	0.00	0.00

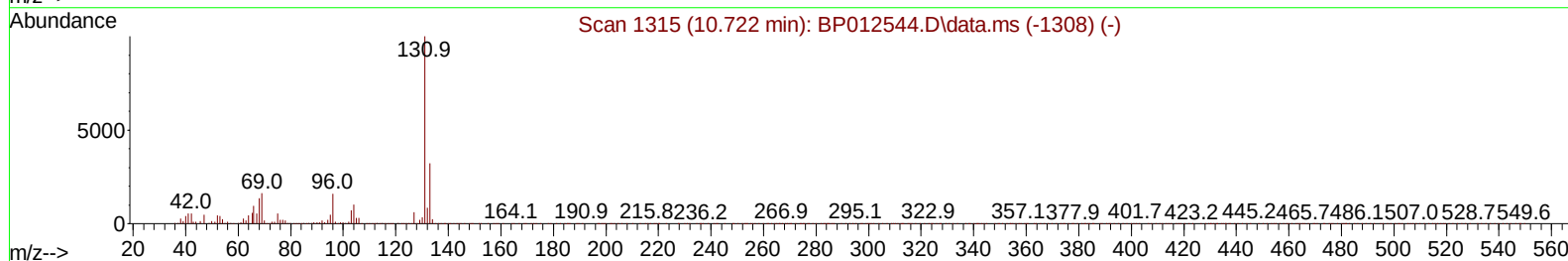
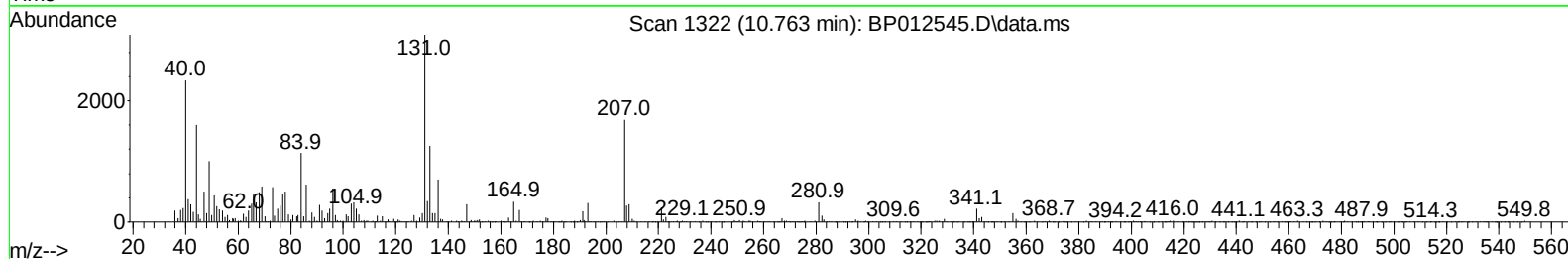
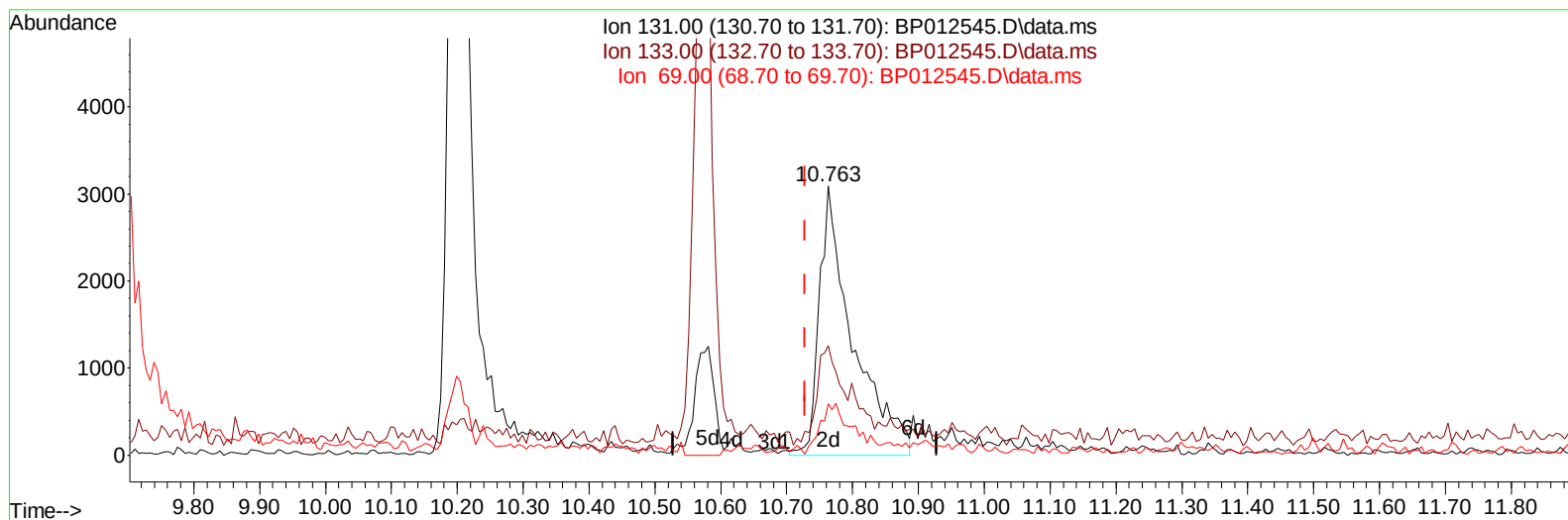
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ALS Vial : 3 Sample Multiplier: 1

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TIC: BP012545.D\data.ms

(31) 4-Chloroaniline-d4 (S)

10.763min (+ 0.035) 0.30 ng/ul m

response 10983

Ion	Exp%	Act%
131.00	100.00	100.00
133.00	32.20	40.73#
69.00	16.50	19.04
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	7.769	152	347558	20.000	ng/ul	0.00
20) Naphthalene-d8	10.575	136	1688490	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.428	164	1170324	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.186	188	2628912	20.000	ng/ul	0.00
79) Chrysene-d12	21.286	240	2291154	20.000	ng/ul	0.00
88) Perylene-d12	23.662	264	2654385	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.228	96	44604	5.167	ng/uL	0.00
4) Pyridine-d5	3.711	84	11452m	0.462	ng/ul	0.06
7) Phenol-d5	6.952	99	831998	26.689	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.110	67	524637	28.192	ng/ul	0.00
11) 2-Chlorophenol-d4	7.304	132	689517	30.289	ng/ul	0.00
15) 4-Methylphenol-d8	8.493	113	694994	28.189	ng/ul	0.00
21) Nitrobenzene-d5	8.946	128	364911	33.327	ng/ul	0.00
24) 2-Nitrophenol-d4	9.663	143	414046	36.432	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.198	165	745862	30.283	ng/ul	0.00
31) 4-Chloroaniline-d4	10.763	131	10983m	0.299	ng/ul	0.04
46) Dimethylphthalate-d6	13.845	166	2562703	31.988	ng/ul	0.00
49) Acenaphthylene-d8	14.116	160	1669550	18.236	ng/ul	0.00
54) 4-Nitrophenol-d4	14.657	143	366375	24.163	ng/ul	0.00
60) Fluorene-d10	15.422	176	2138326	31.174	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.563	200	301507	22.028	ng/ul	0.00
73) Anthracene-d10	17.286	188	3152396	27.830	ng/ul	0.00
81) Pyrene-d10	19.533	212	3696920	32.138	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.510	264	3350974	25.837	ng/ul	0.00

Target Compounds Qvalue

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

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