

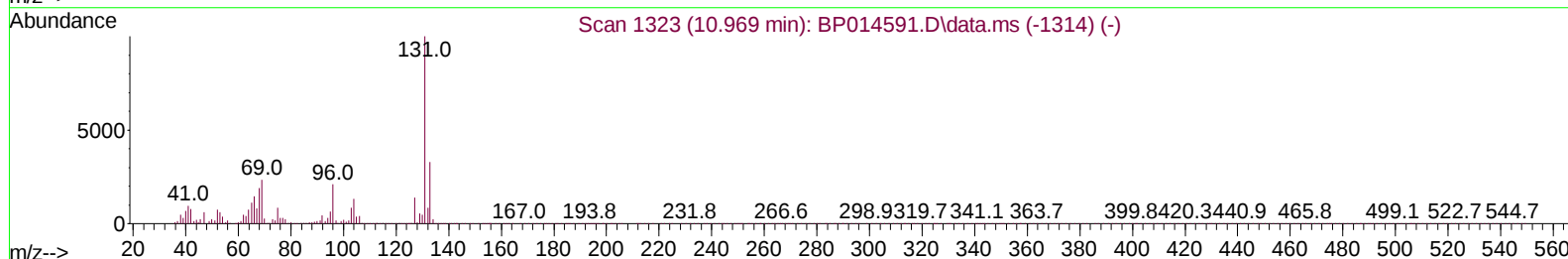
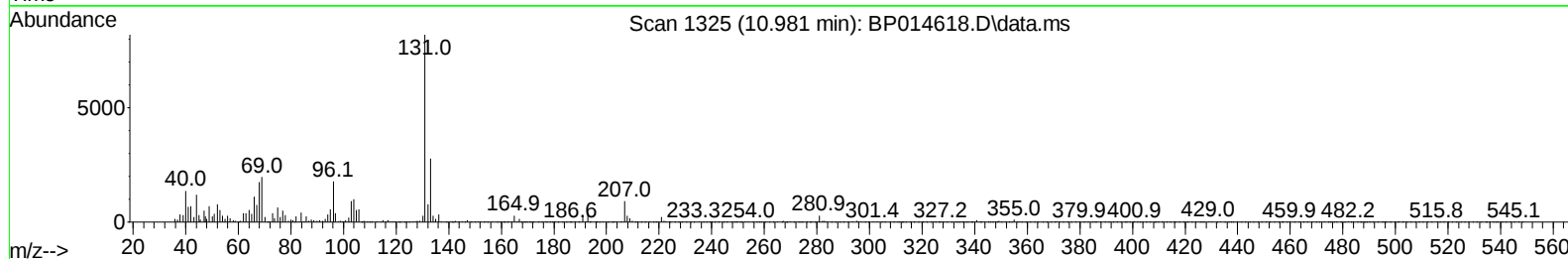
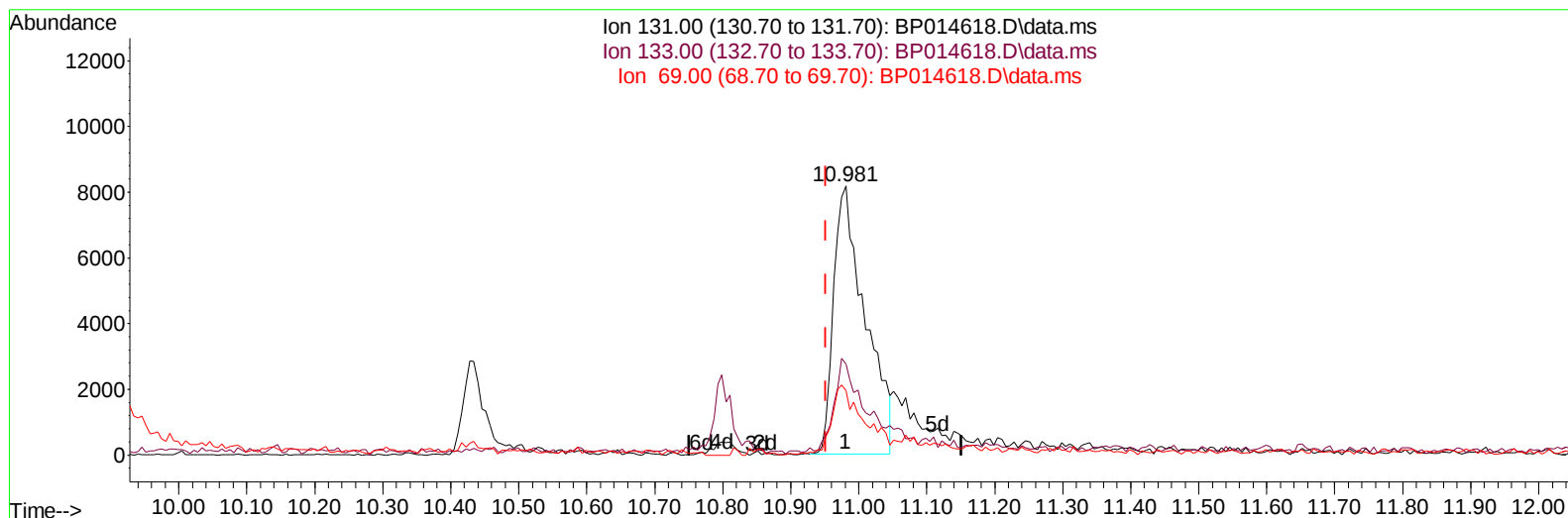
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP040723\
 Data File : BP014618.D
 Acq On : 07 Apr 2023 22:55
 Operator : CG/JU
 Sample : 02225-12
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0AA9

Manual IntegrationsAPPROVED

Quant Time: Apr 08 01:51:45 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP032823.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Apr 07 00:45:34 2023
 Response via : Initial Calibration

Reviewed By :Christian Giraldo 04/10/2023
 Supervised By :Jagrut Upadhyay 04/10/2023



TIC: BP014618.D\data.ms

(31) 4-Chloroaniline-d4 (S)

10.981min (+ 0.029) 2.41 ng/ul

response 26320

Ion	Exp%	Act%
131.00	100.00	100.00
133.00	32.20	33.76
69.00	21.30	23.92
0.00	0.00	0.00

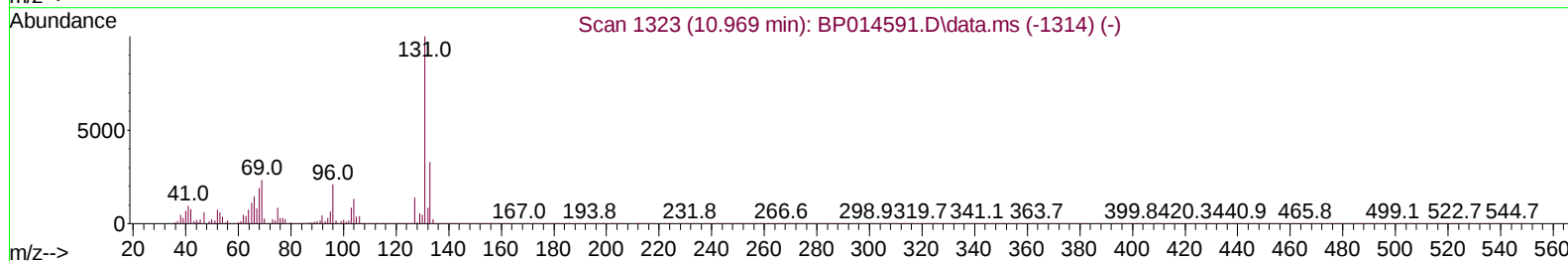
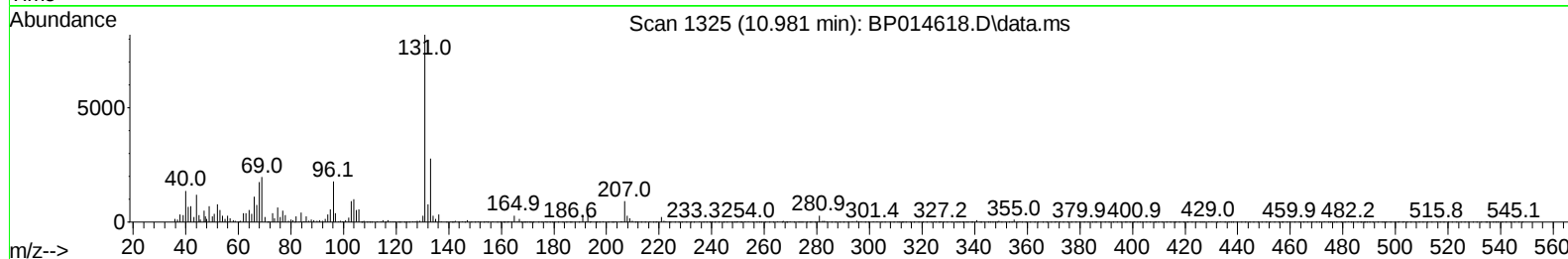
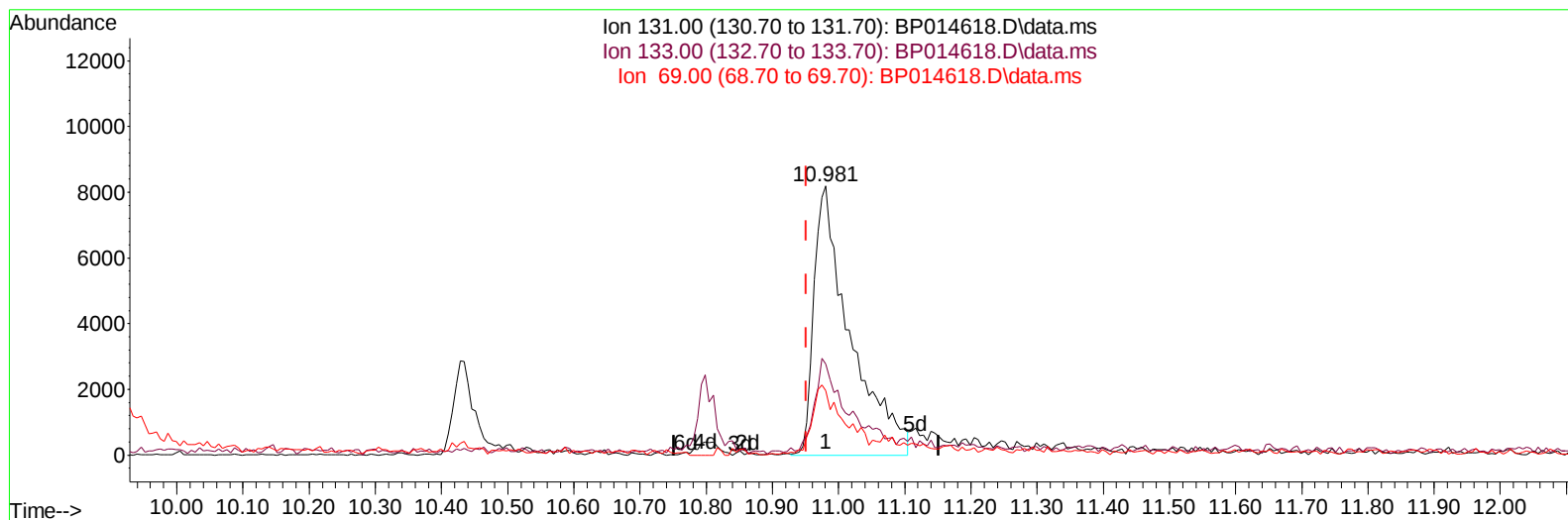
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TIC: BP014618.D\data.ms

(31) 4-Chloroaniline-d4 (S)

10.981min (+ 0.029) 2.84 ng/ul m

response 31024

Ion	Exp%	Act%
131.00	100.00	100.00
133.00	32.20	33.76
69.00	21.30	23.92
0.00	0.00	0.00

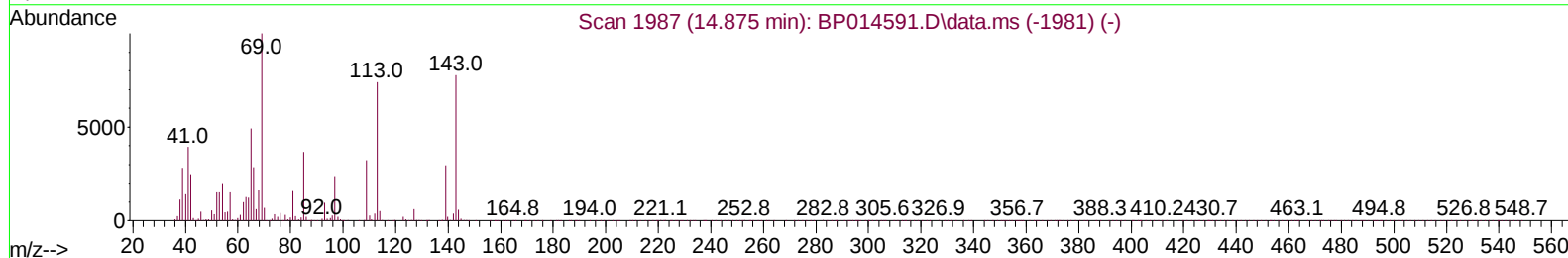
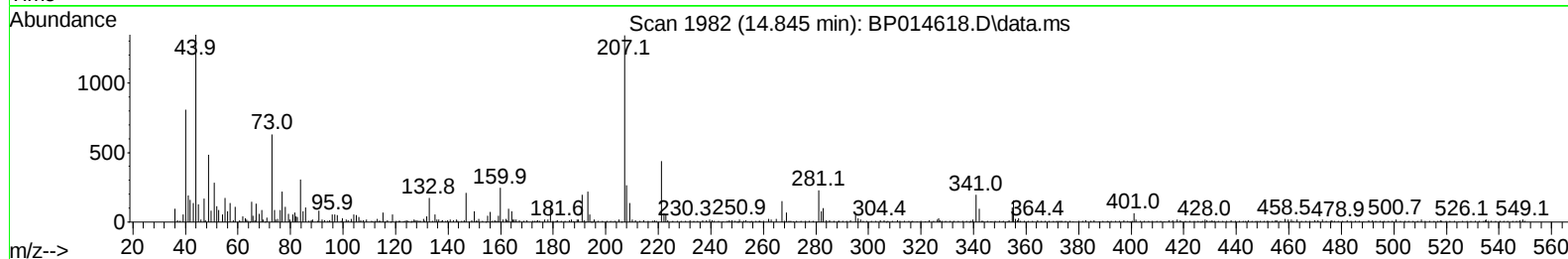
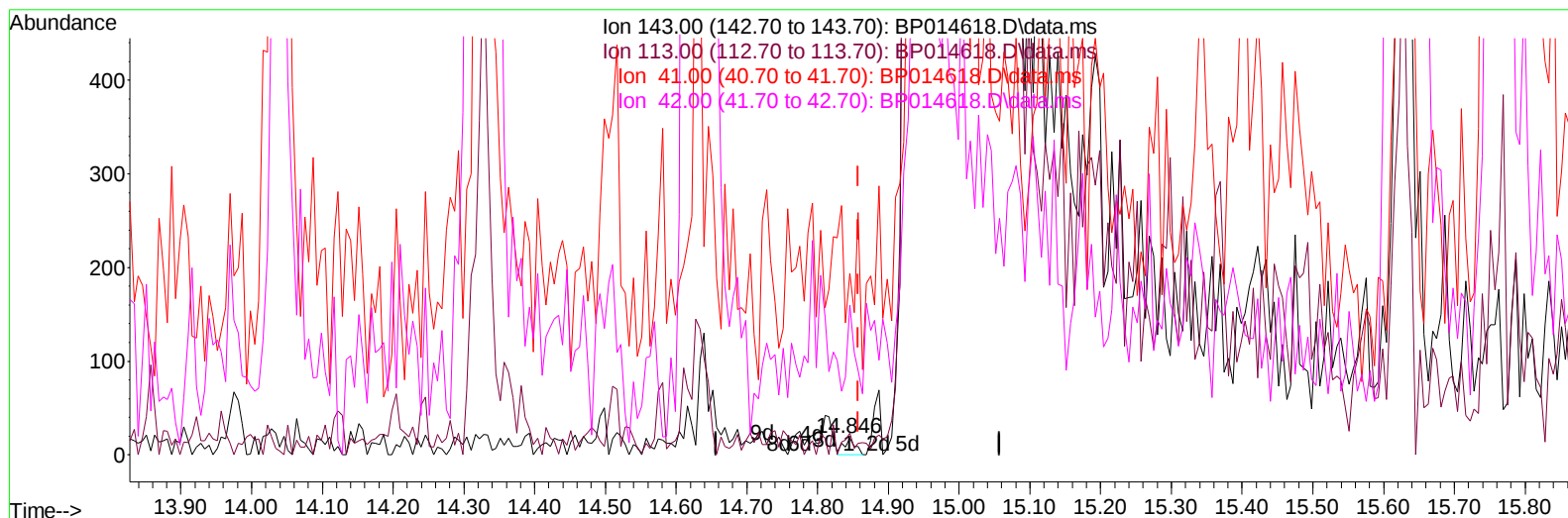
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Instrument :
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TIC: BP014618.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.845min (-0.012) 0.01 ng/u1

response 20

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	96.80	111.11
41.00	44.20	718.52#
42.00	31.50	592.59#

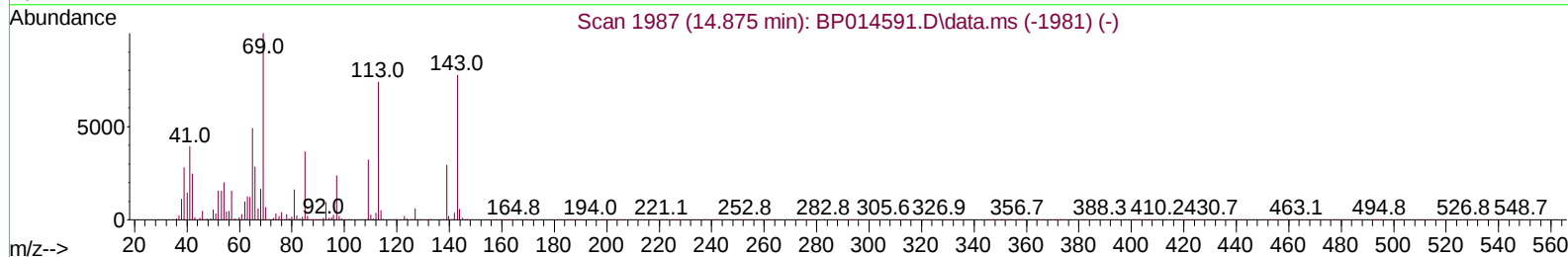
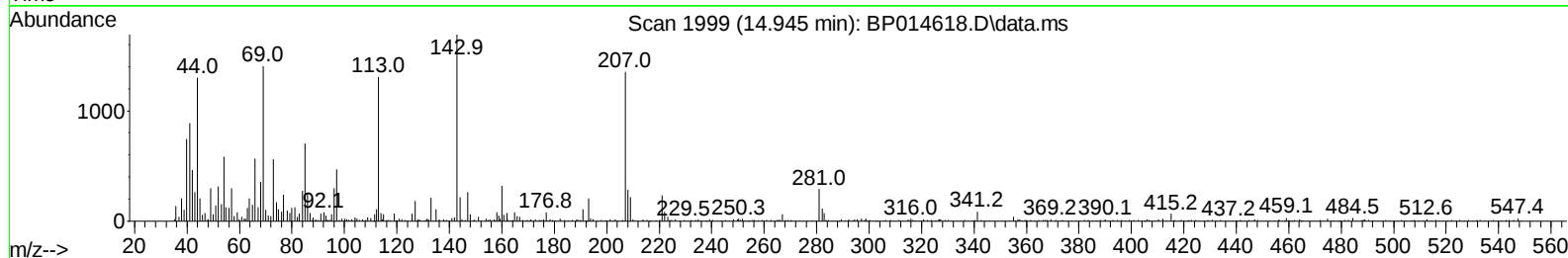
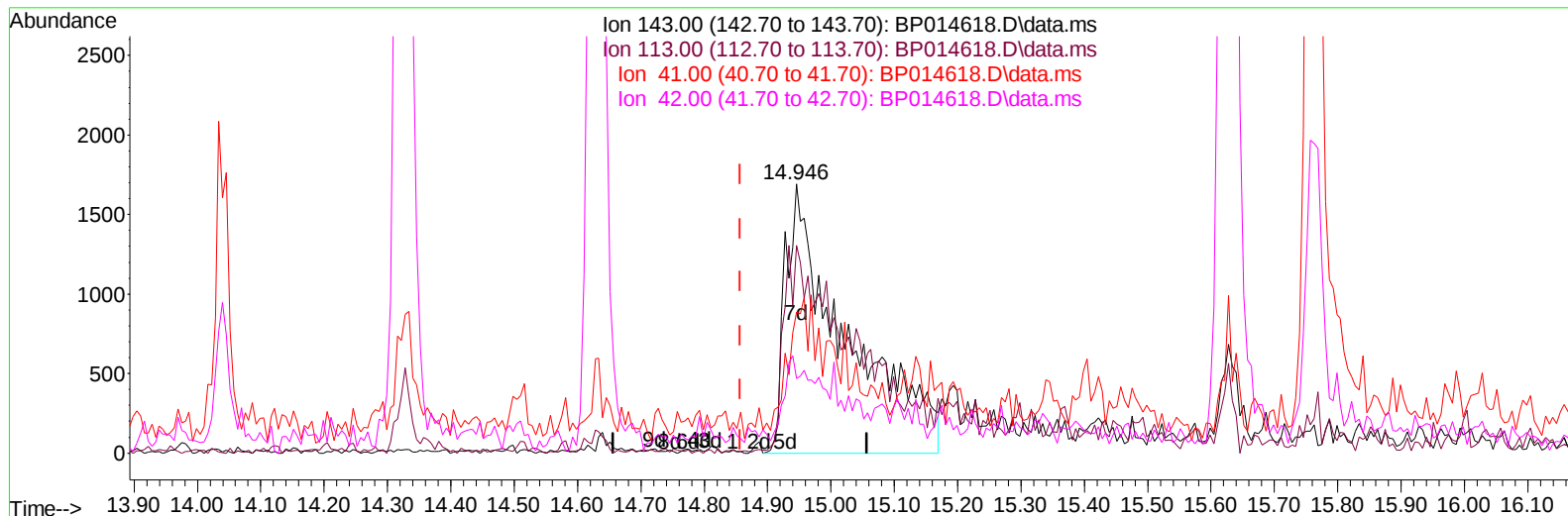
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Sample : 02225-12
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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TIC: BP014618.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.945min (+ 0.088) 2.88 ng/ul m

response 11157

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	96.80	77.23#
41.00	44.20	52.45
42.00	31.50	27.62

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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	7.981	152	116064	20.000	ng/ul	0.00
20) Naphthalene-d8	10.799	136	492017	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.634	164	316670	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.392	188	673867	20.000	ng/ul	0.00
79) Chrysene-d12	21.480	240	619928	20.000	ng/ul	0.00
88) Perylene-d12	23.986	264	749160	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.364	96	12676	4.235	ng/uL	0.00
4) Pyridine-d5	0.000	84	0d	0.000	ng/ul	
7) Phenol-d5	7.158	99	45637	4.264	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.305	67	207755	30.251	ng/ul	0.00
11) 2-Chlorophenol-d4	7.511	132	161159	20.831	ng/ul	0.00
15) 4-Methylphenol-d8	8.705	113	96976	11.232	ng/ul	0.00
21) Nitrobenzene-d5	9.158	128	106650	26.851	ng/ul	0.00
24) 2-Nitrophenol-d4	9.881	143	116718	25.662	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.428	165	178861	21.593	ng/ul	0.00
31) 4-Chloroaniline-d4	10.981	131	31024m	2.843	ng/ul	0.03
46) Dimethylphthalate-d6	14.040	166	752887	29.322	ng/ul	0.00
49) Acenaphthylene-d8	14.328	160	824155	28.927	ng/ul	0.00
54) 4-Nitrophenol-d4	14.945	143	11157m	2.876	ng/ul	0.09
60) Fluorene-d10	15.628	176	631638	29.499	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.763	200	111062	22.399	ng/ul	0.00
73) Anthracene-d10	17.492	188	1053111	32.292	ng/ul	0.00
81) Pyrene-d10	19.722	212	1265601	30.394	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.822	264	1310250	32.919	ng/ul	0.00

Target Compounds Qvalue

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

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