

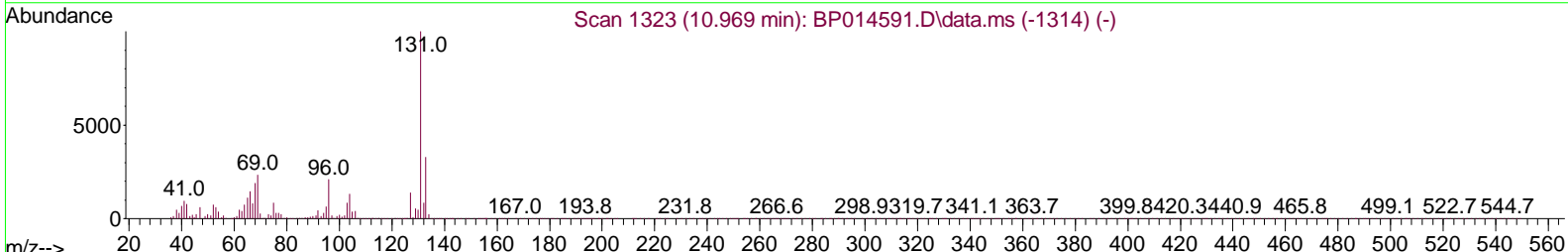
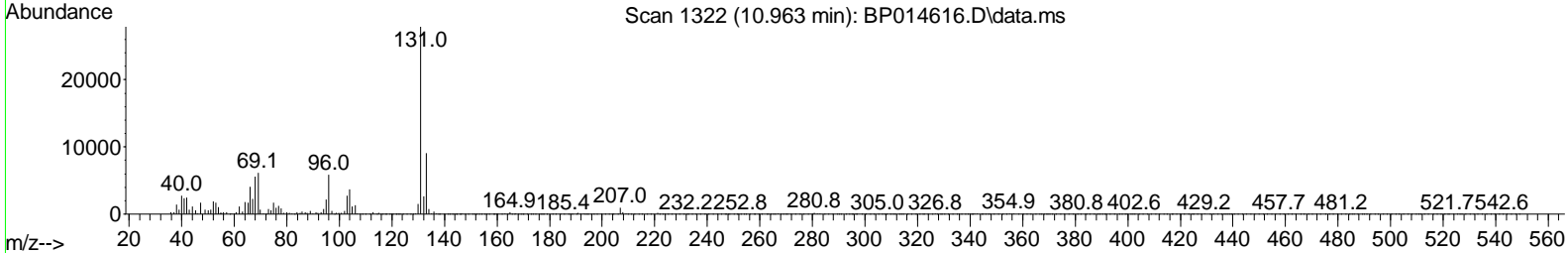
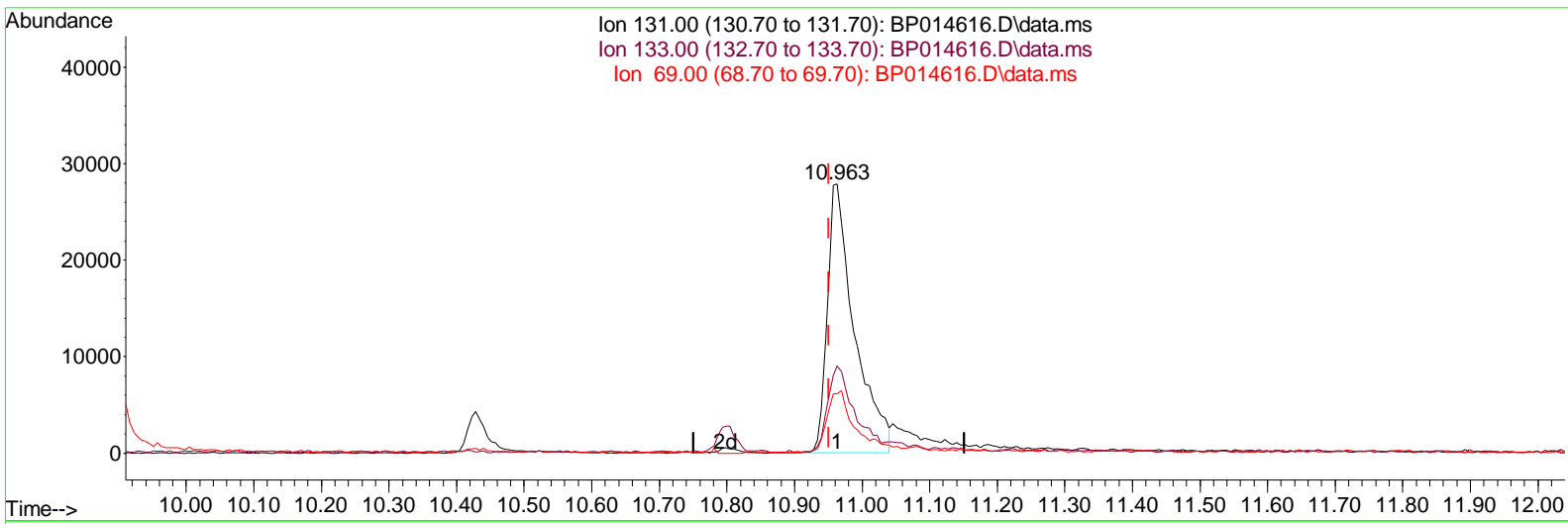
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP040723\
 Data File : BP014616.D
 Acq On : 07 Apr 2023 21:47
 Operator : CG/JU
 Sample : 02225-21
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0AE6

Manual Integrations APPROVED

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

Quant Time: Apr 08 01:51:17 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



TIC: BP014616.D\data.ms

(31) 4-Chloroaniline-d4 (S)

10.963min (+ 0.012) 5.62 ng/ul

response 77469

Ion	Exp%	Act%
131.00	100.00	100.00
133.00	32.20	32.53
69.00	21.30	22.10
0.00	0.00	0.00

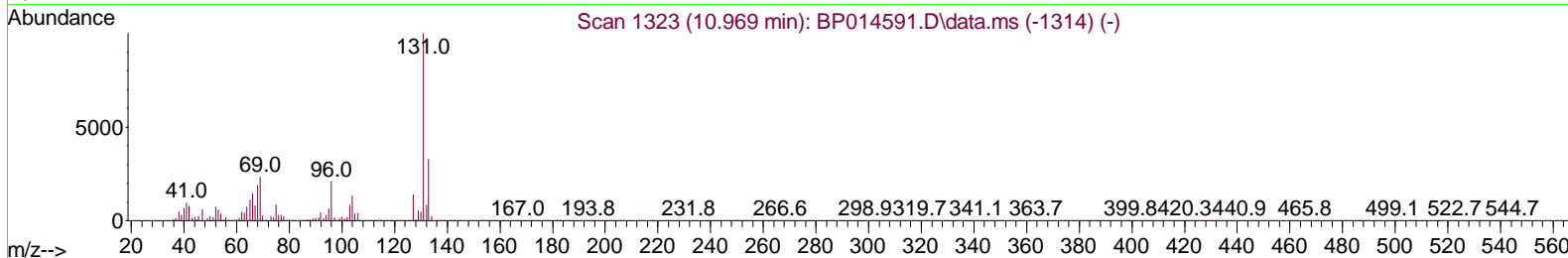
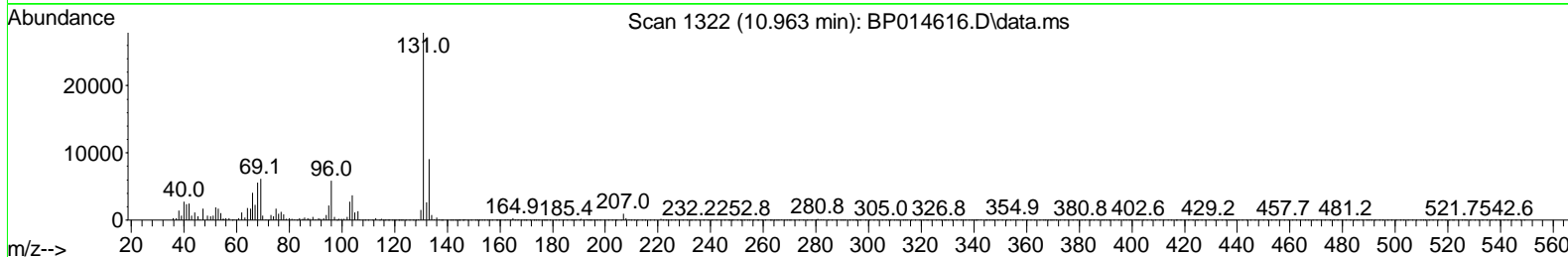
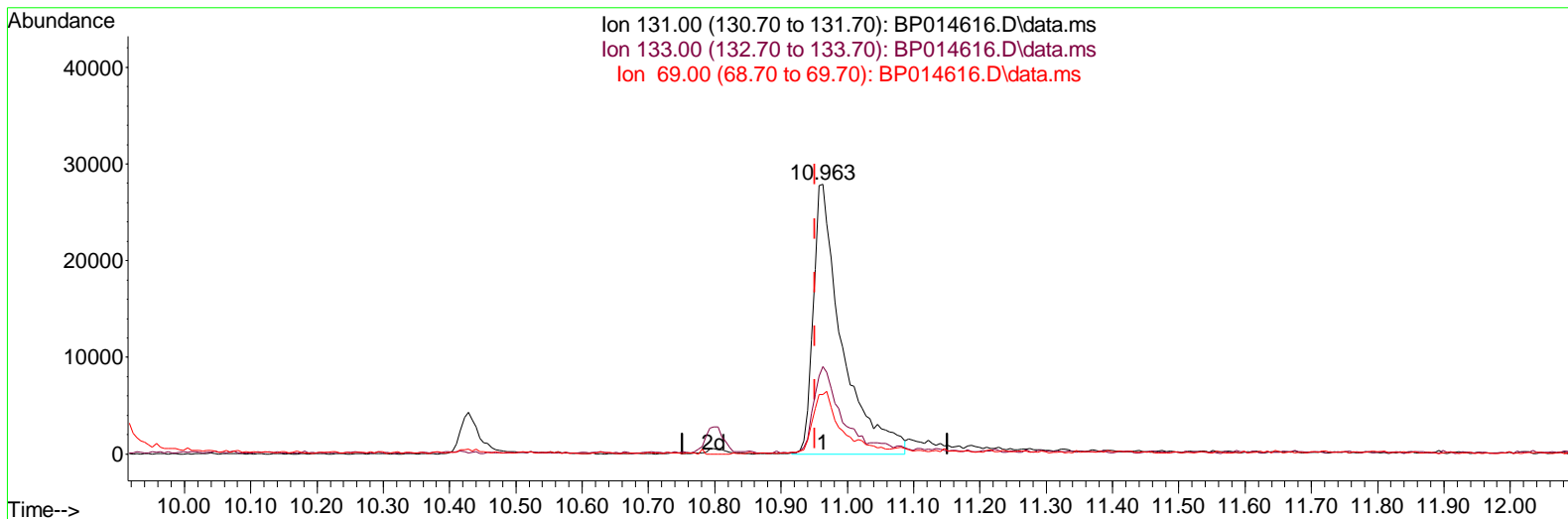
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TIC: BP014616.D\data.ms

(31) 4-Chloroaniline-d4 (S)

10.963min (+ 0.012) 6.08 ng/ul m

response 83791

Ion	Exp%	Act%
131.00	100.00	100.00
133.00	32.20	32.53
69.00	21.30	22.10
0.00	0.00	0.00

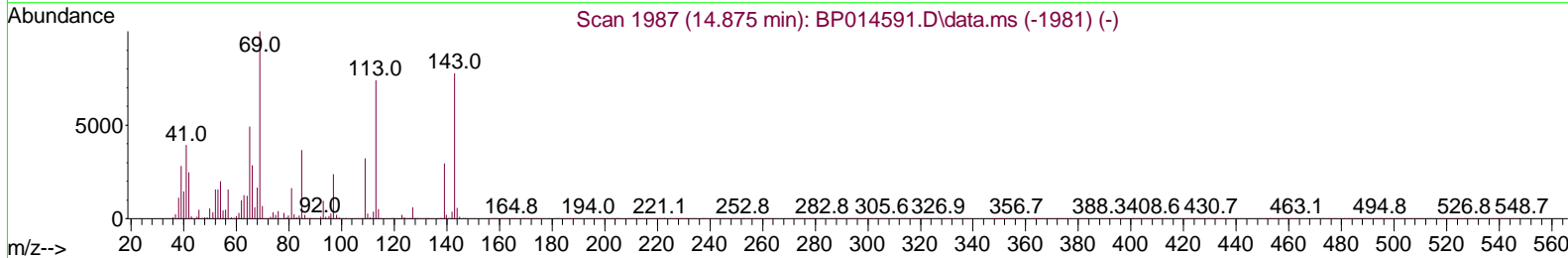
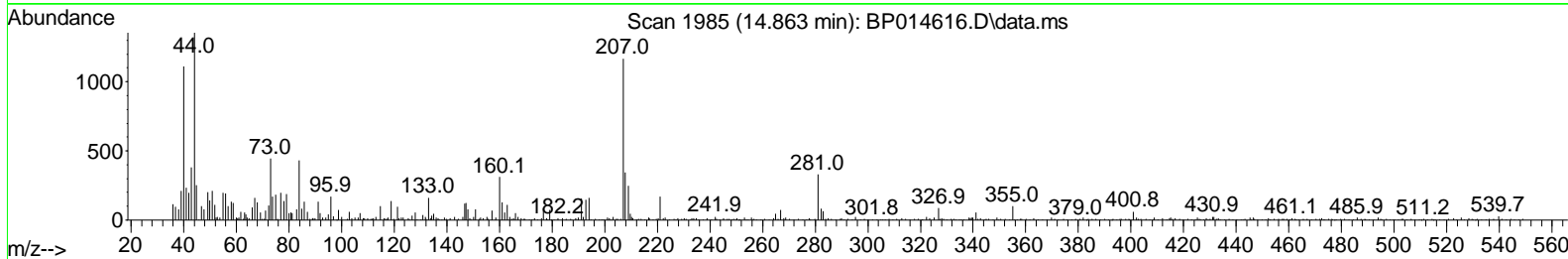
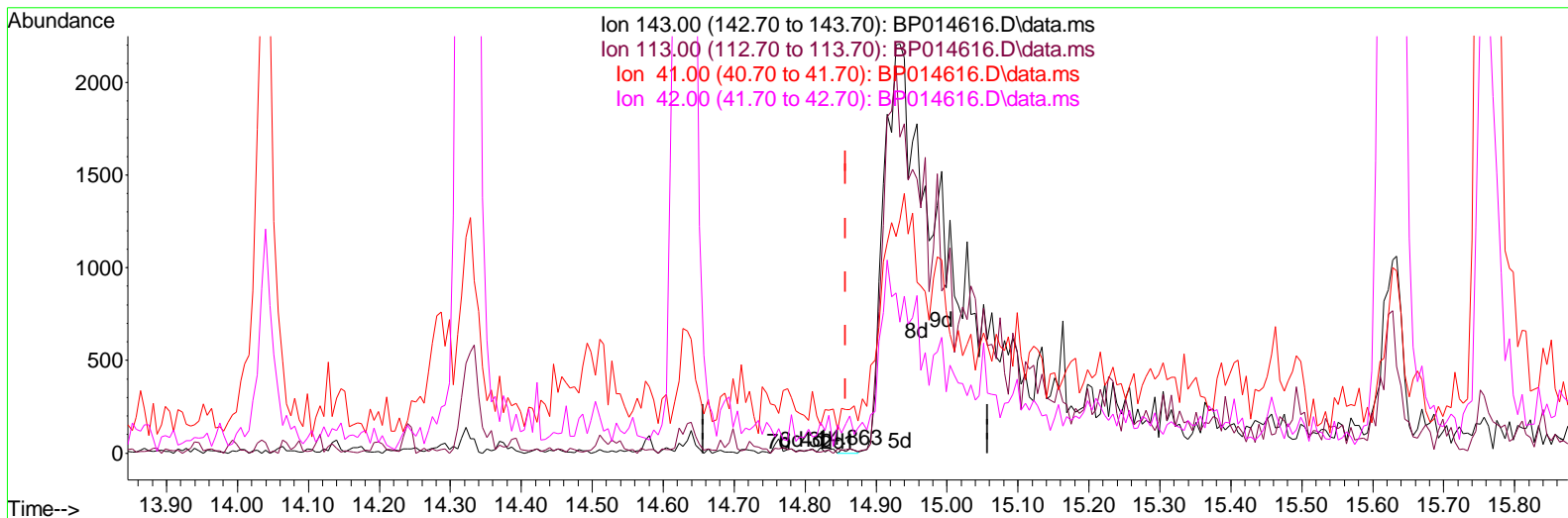
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 ALS Vial : 9 Sample Multiplier: 1

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TIC: BP014616.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.863min (+ 0.006) 0.00 ng/ul

response 25

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	96.80	114.29
41.00	44.20	1109.52#
42.00	31.50	947.62#

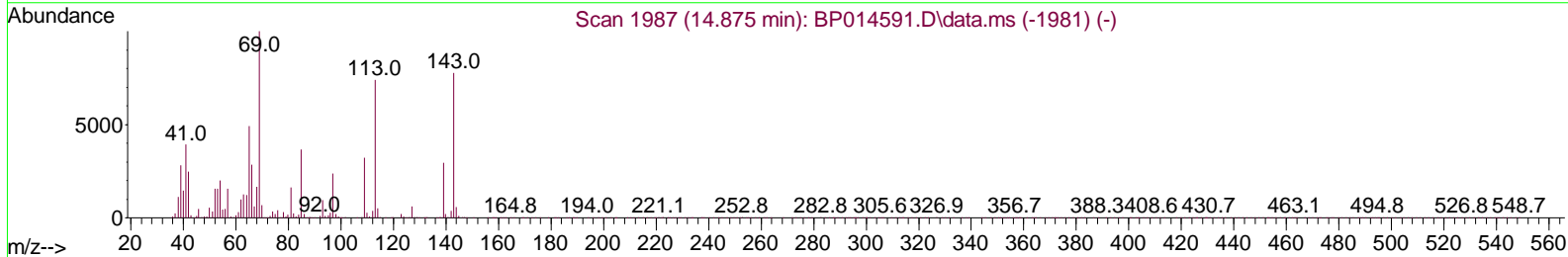
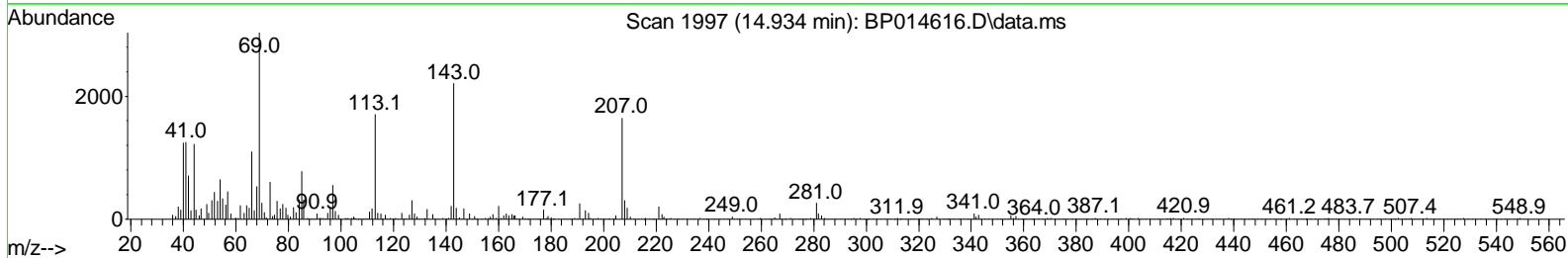
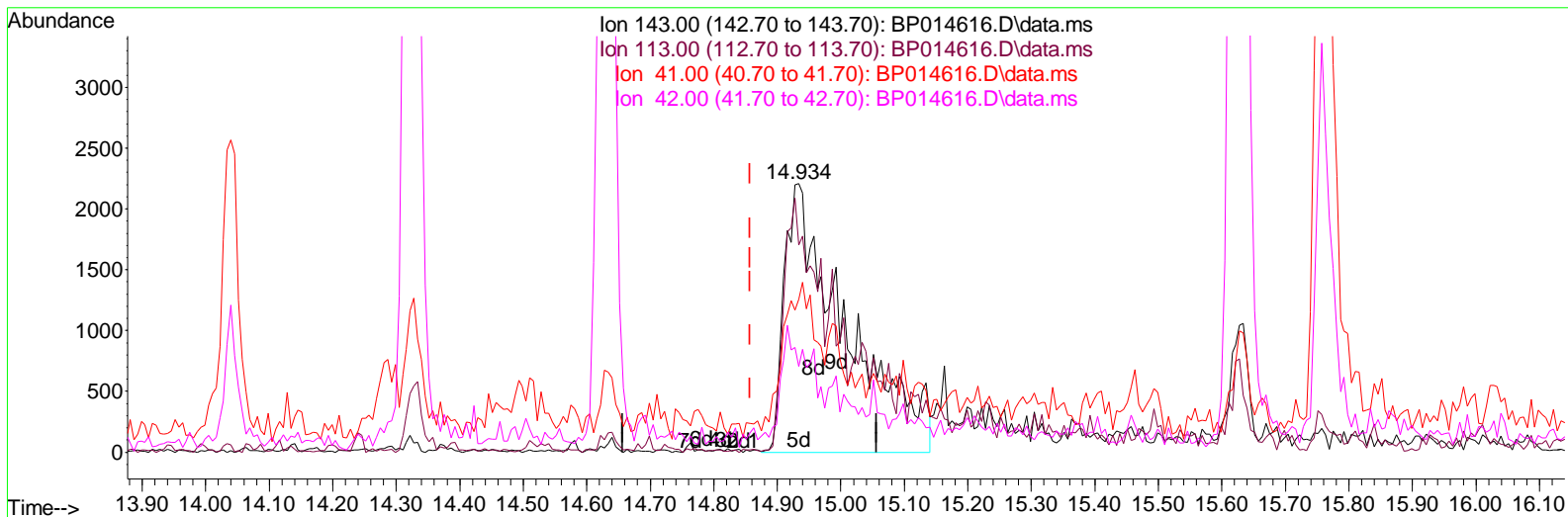
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 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
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TIC: BP014616.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.934min (+ 0.077) 2.94 ng/ul m

response	14721	
Ion	Exp%	Act%
143.00	100.00	100.00
113.00	96.80	77.16#
41.00	44.20	57.00#
42.00	31.50	31.99

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Compound	R. T.	QI on	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.981	152	141054	20.000	ng/ul	0.00
20) Naphthalene-d8	10.799	136	621621	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.634	164	408285	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.392	188	863797	20.000	ng/ul	0.00
79) Chrysene-d12	21.486	240	651223	20.000	ng/ul	0.00
88) Perylene-d12	23.986	264	722790	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.364	96	16026	4.406	ng/uL	0.00
4) Pyridine-d5	0.000	84	0d	0.000	ng/ul	
7) Phenol-d5	7.158	99	55726	4.284	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.305	67	249285	29.867	ng/ul	0.00
11) 2-Chlorophenol-d4	7.511	132	197046	20.957	ng/ul	0.00
15) 4-Methylphenol-d8	8.705	113	119182	11.358	ng/ul	0.00
21) Nitrobenzene-d5	9.158	128	132131	26.330	ng/ul	0.00
24) 2-Nitrophenol-d4	9.881	143	146893	25.563	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.428	165	233236	22.287	ng/ul	0.00
31) 4-Chloroaniline-d4	10.963	131	83791m	6.079	ng/ul	0.01
46) Dimethylphthalate-d6	14.040	166	1012307	30.579	ng/ul	0.00
49) Acenaphthylene-d8	14.328	160	1094051	29.784	ng/ul	0.00
54) 4-Nitrophenol-d4	14.934	143	14721m	2.944	ng/ul	0.08
60) Fluorene-d10	15.628	176	835626	30.269	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.763	200	151295	23.804	ng/ul	0.00
73) Anthracene-d10	17.492	188	1320552	31.589	ng/ul	0.00
81) Pyrene-d10	19.722	212	1455123	33.266	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.822	264	1258798	32.780	ng/ul	0.00

Target Compounds Qvalue

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

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