

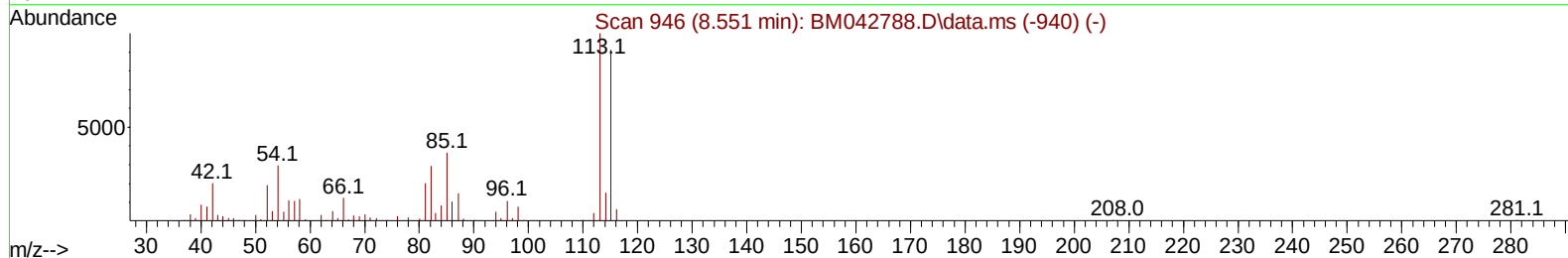
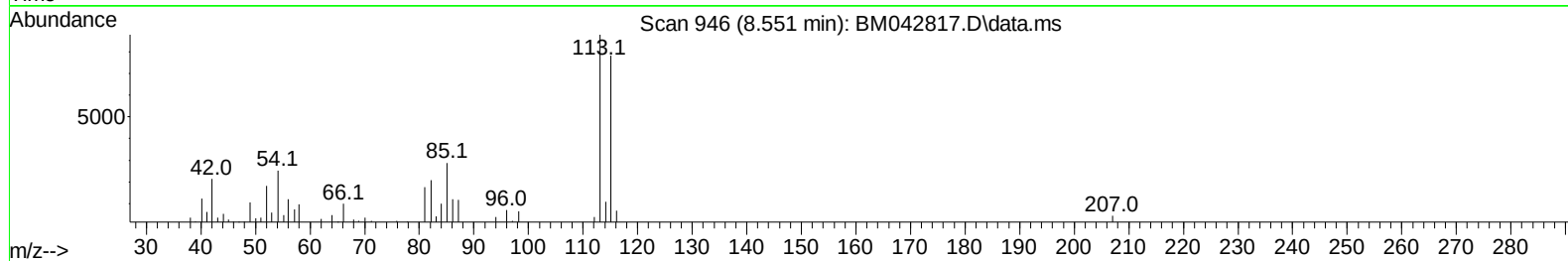
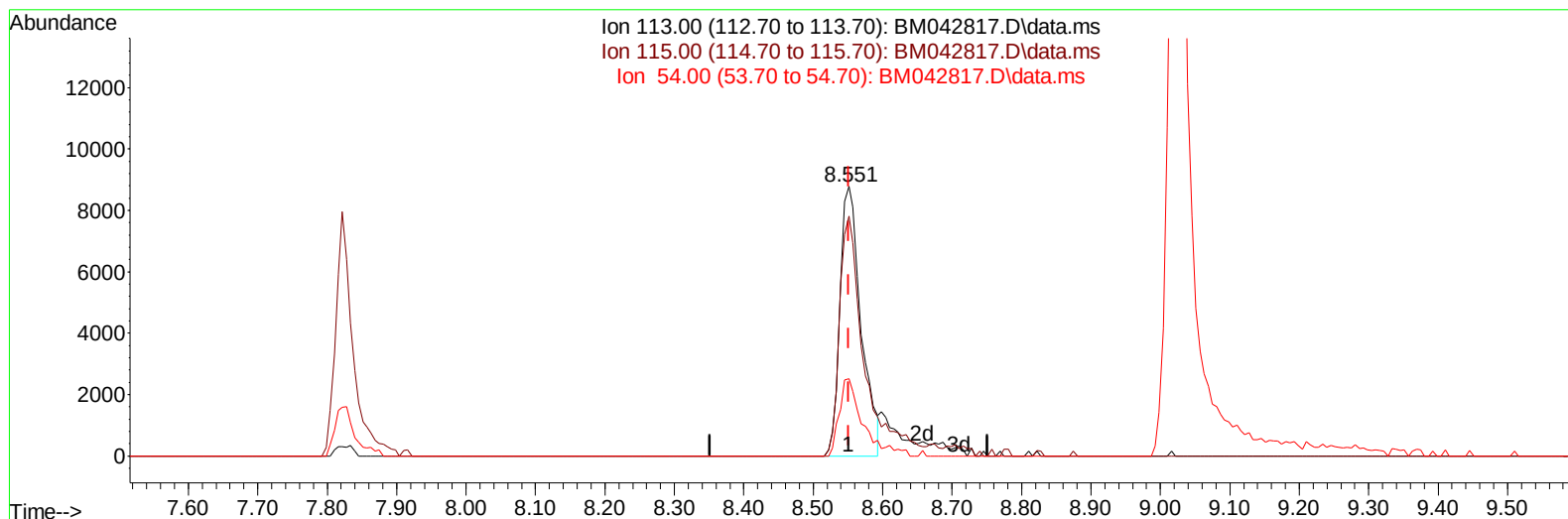
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM111323\
 Data File : BM042817.D
 Acq On : 16 Nov 2023 22:08
 Operator : MA/JU
 Sample : 05268-08
 Misc :
 ALS Vial : 59 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 DCKD0

Manual IntegrationsAPPROVED

Quant Time: Nov 16 23:56:32 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM111323.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Nov 15 22:28:07 2023
 Response via : Initial Calibration

Reviewed By :Yogesh Patel 11/17/2023
 Supervised By :mohammad ahmed 11/19/2023



TIC: BM042817.D\data.ms

(15) 4-Methylphenol-d8 (S)

8.551min (-0.000) 11.70 ng/u1

response 18402

Ion	Exp%	Act%
113.00	100.00	100.00
115.00	91.70	89.07
54.00	24.00	28.58
0.00	0.00	0.00

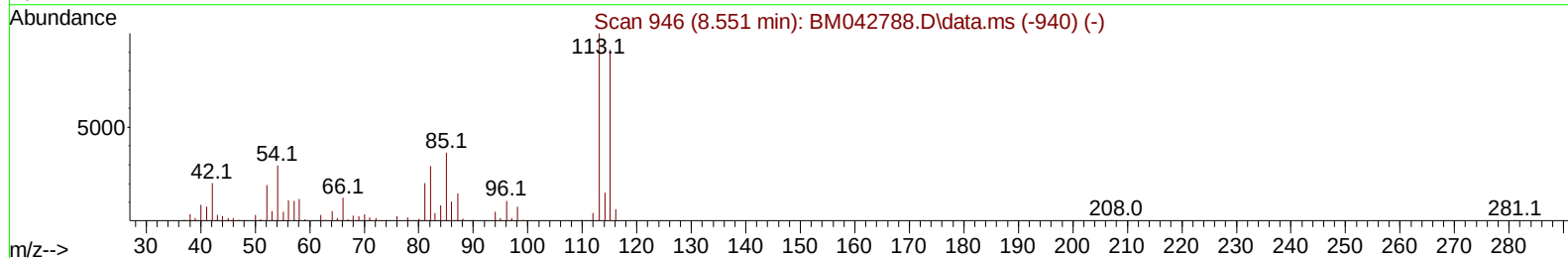
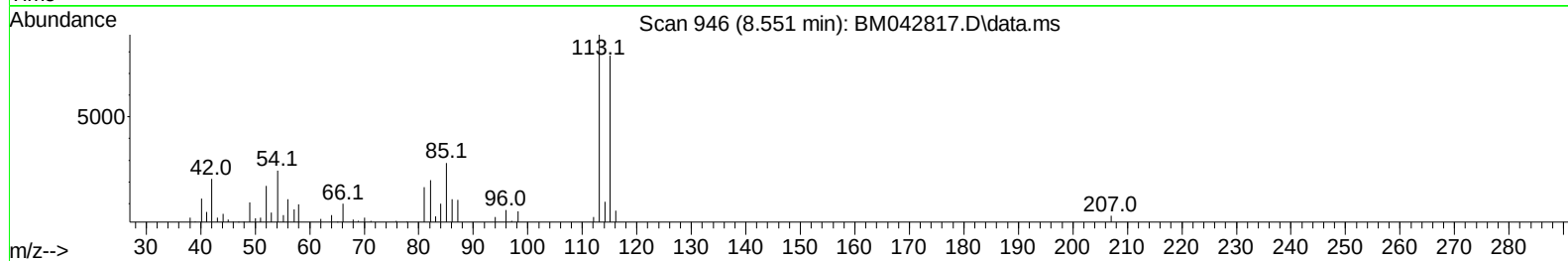
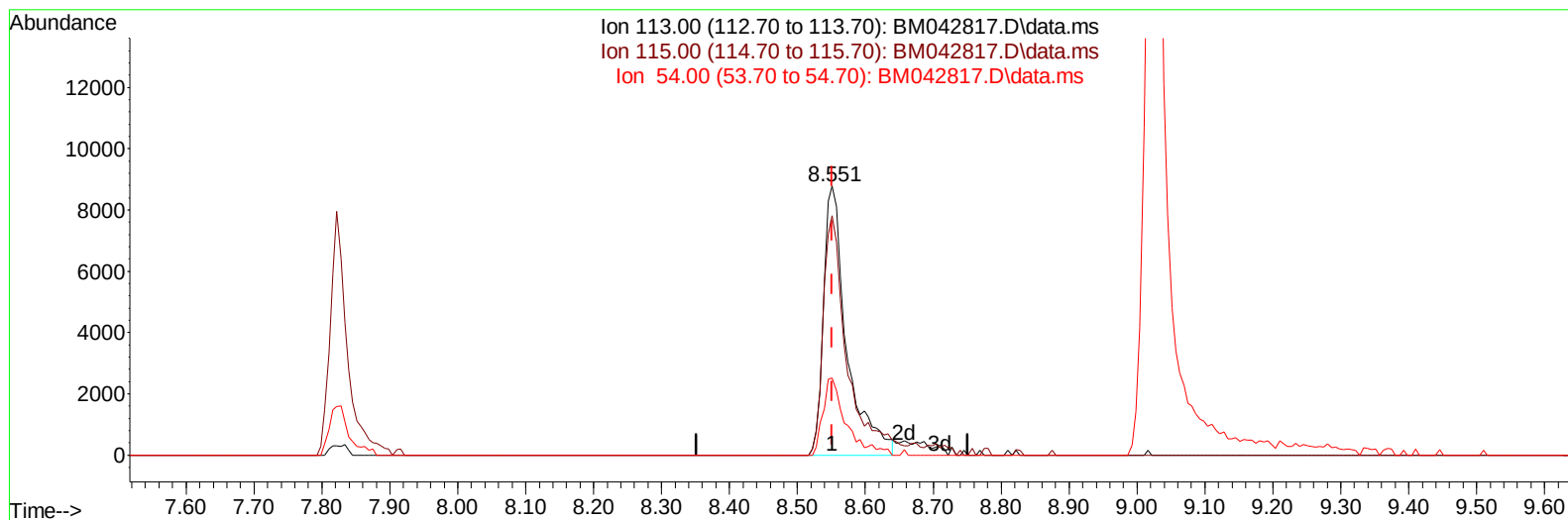
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TIC: BM042817.D\data.ms

(15) 4-Methylphenol-d8 (S)

8.551min (-0.000) 13.23 ng/ul m

response 20811

Ion	Exp%	Act%
113.00	100.00	100.00
115.00	91.70	89.07
54.00	24.00	28.58
0.00	0.00	0.00

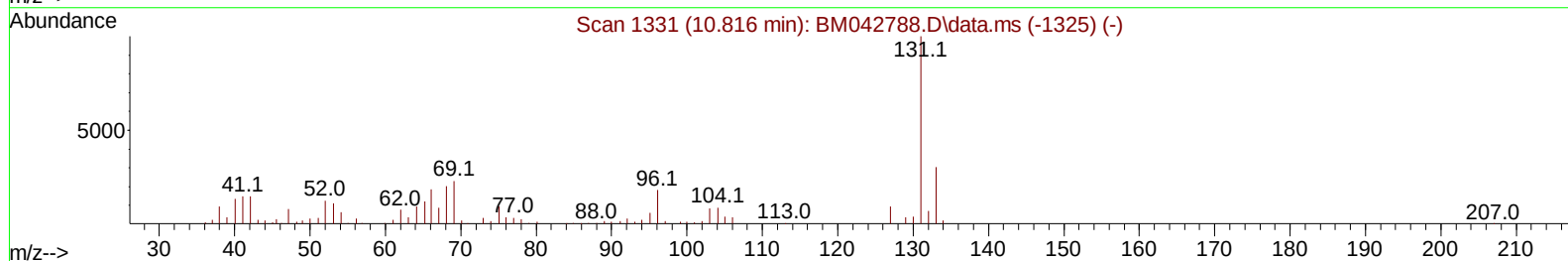
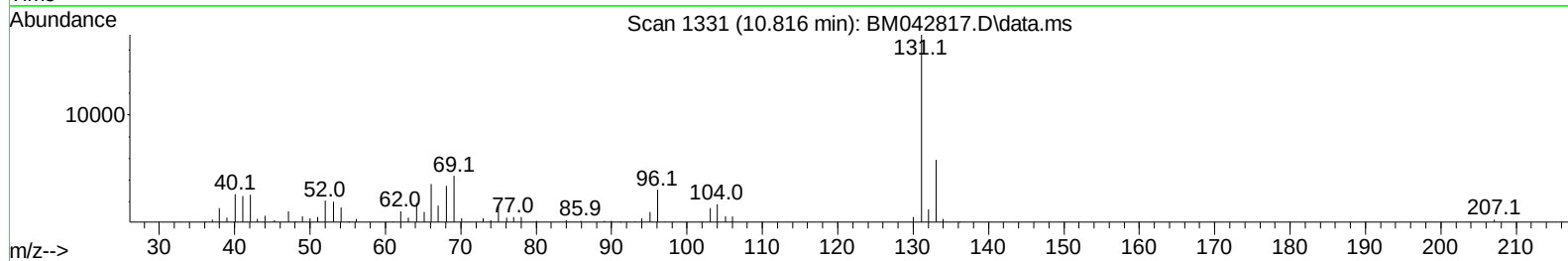
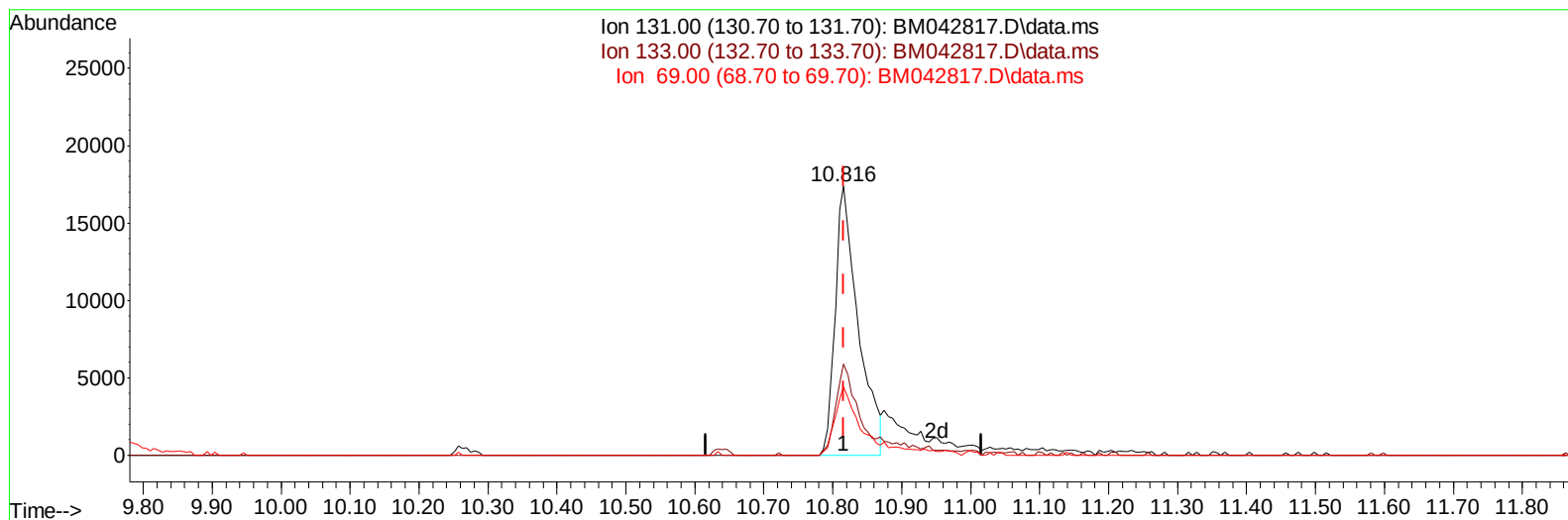
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 ALS Vial : 59 Sample Multiplier: 1

Instrument :
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TIC: BM042817.D\data.ms

(31) 4-Chloroaniline-d4 (S)

10.816min (-0.000) 19.72 ng/u1

response 40215

Ion	Exp%	Act%
131.00	100.00	100.00
133.00	33.00	33.86
69.00	23.40	25.39
0.00	0.00	0.00

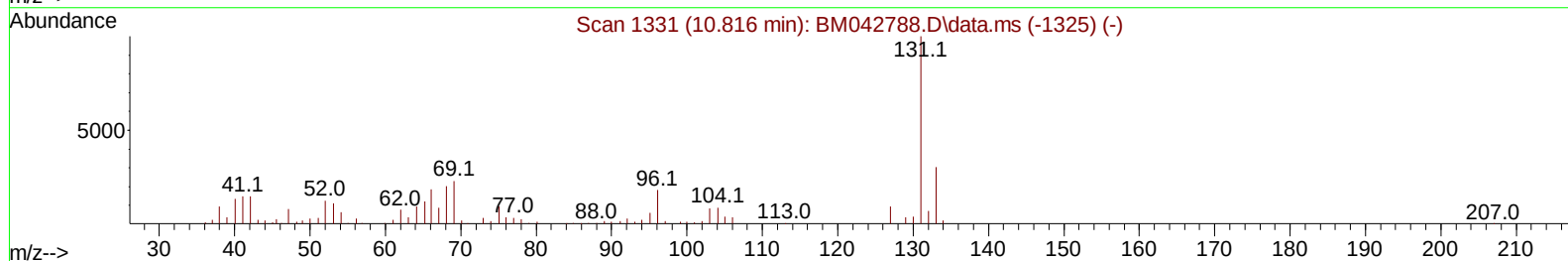
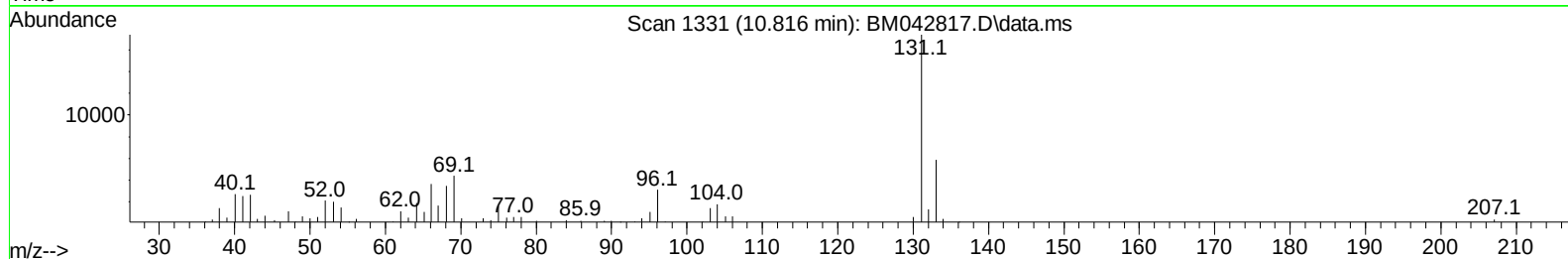
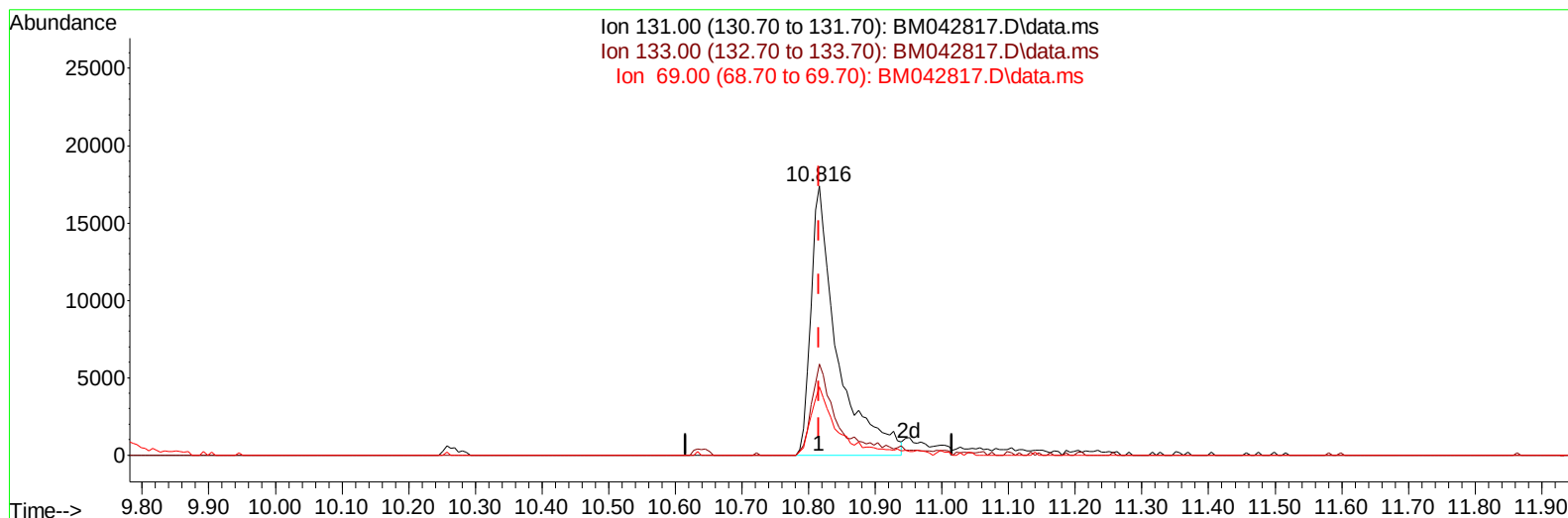
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 Misc :
 ALS Vial : 59 Sample Multiplier: 1

Instrument :
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Manual IntegrationsAPPROVED

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TIC: BM042817.D\data.ms

(31) 4-Chloroaniline-d4 (S)

10.816min (-0.000) 23.32 ng/ul m

response 47563

Ion	Exp%	Act%
131.00	100.00	100.00
133.00	33.00	33.86
69.00	23.40	25.39
0.00	0.00	0.00

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Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 11/17/2023
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Quant Time: Nov 17 00:41:45 2023
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.822	152	20521	20.000	ng/ul	0.00
20) Naphthalene-d8	10.634	136	84220	20.000	ng/ul	# 0.00
38) Acenaphthene-d10	14.480	164	53427	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.233	188	118576	20.000	ng/ul	0.00
79) Chrysene-d12	21.415	240	118953	20.000	ng/ul	0.00
88) Perylene-d12	23.774	264	144993	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.175	96	1828	2.795	ng/uL	0.00
4) Pyridine-d5	3.622	84	14537	8.936	ng/ul	0.00
7) Phenol-d5	7.004	99	9151	4.582	ng/ul	0.01
9) Bis-(2-Chloroethyl)eth...	7.169	67	52105	37.128	ng/ul	0.00
11) 2-Chlorophenol-d4	7.345	132	37094	24.070	ng/ul	0.00
15) 4-Methylphenol-d8	8.551	113	20811m	13.234	ng/ul	0.00
21) Nitrobenzene-d5	9.028	128	23318	31.579	ng/ul	0.00
24) 2-Nitrophenol-d4	9.739	143	25071	29.024	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.263	165	39213	24.414	ng/ul	0.00
31) 4-Chloroaniline-d4	10.816	131	47563m	23.323	ng/ul	0.00
46) Dimethylphthalate-d6	13.904	166	189820	37.540	ng/ul	0.00
49) Acenaphthylene-d8	14.180	160	193066	35.685	ng/ul	0.00
54) 4-Nitrophenol-d4	14.798	143	114	0.200	ng/ul	0.06
60) Fluorene-d10	15.474	176	164711	39.337	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.621	200	21785	24.255	ng/ul	0.00
73) Anthracene-d10	17.333	188	272599	41.021	ng/ul	0.00
81) Pyrene-d10	19.627	212	367650	47.198	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.621	264	389302	44.887	ng/ul	0.00

Target Compounds Qvalue

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

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