

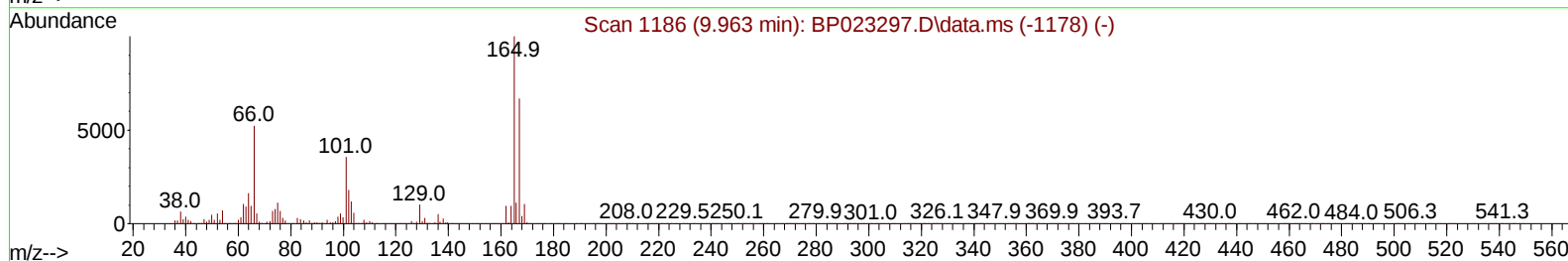
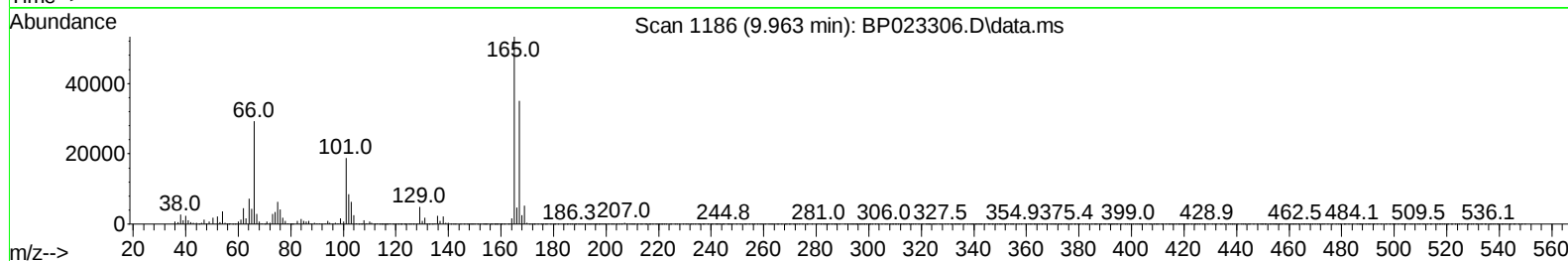
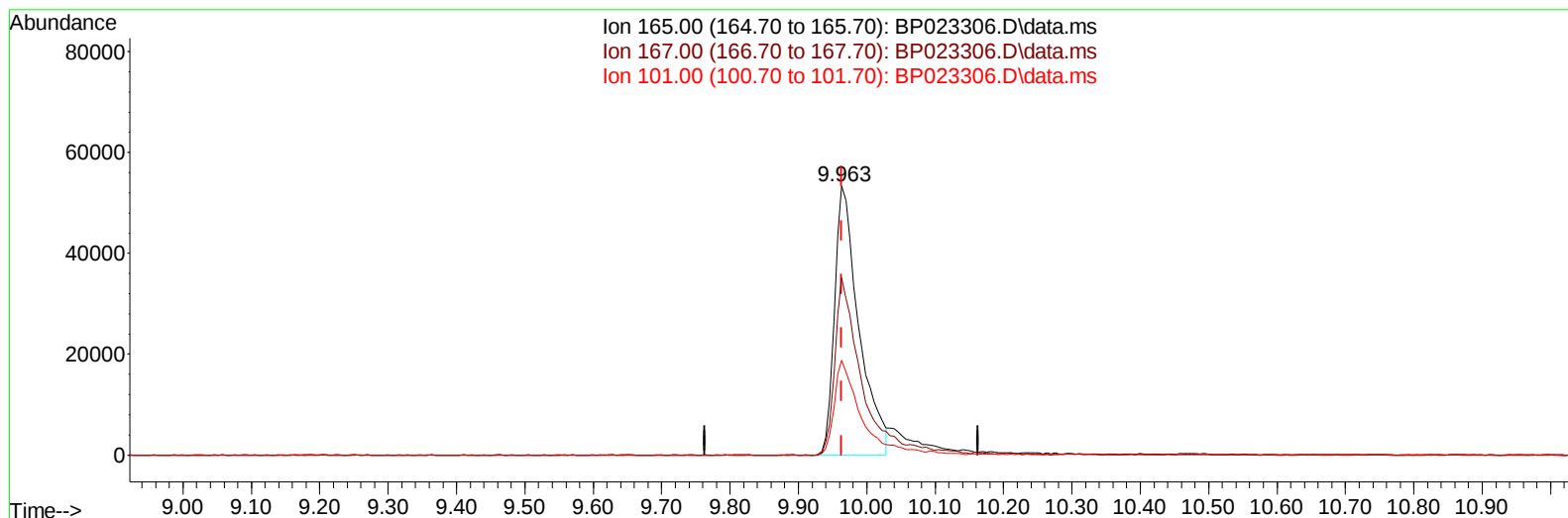
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP112224\
Data File : BP023306.D
Acq On : 22 Nov 2024 16:43
Operator : RC/JU
Sample : PB165140BL
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
SBLK140

Manual IntegrationsAPPROVED

Quant Time: Nov 22 23:46:43 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP110924.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed Nov 20 04:34:35 2024
Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 11/25/2024
Supervised By :mohammad ahmed 11/27/2024



TIC: BP023306.D\data.ms

(28) 2,4-Dichlorophenol-d3 (S)

9.963min (+ 0.000) 29.60 ng/u1

response 131778

Ion	Exp%	Act%
165.00	100.00	100.00
167.00	65.80	65.90
101.00	32.60	35.17
0.00	0.00	0.00

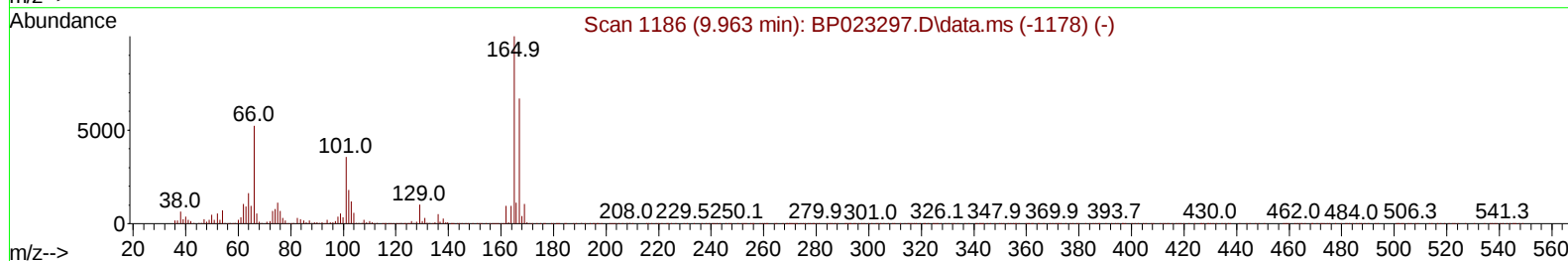
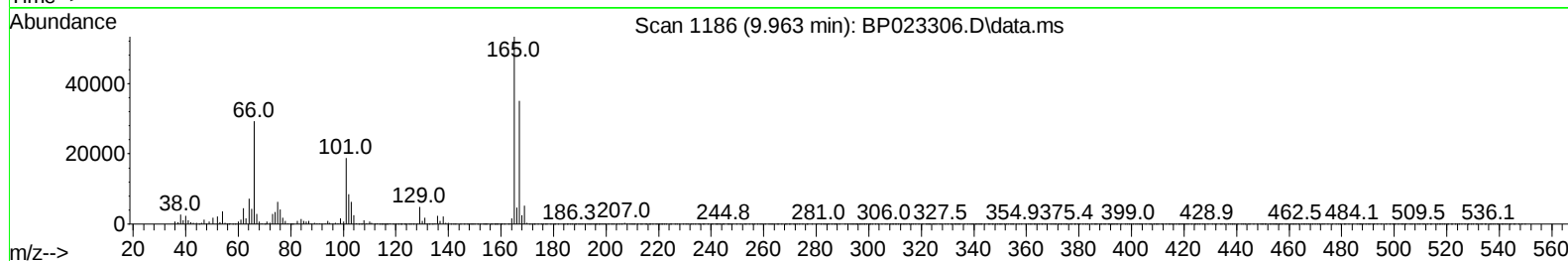
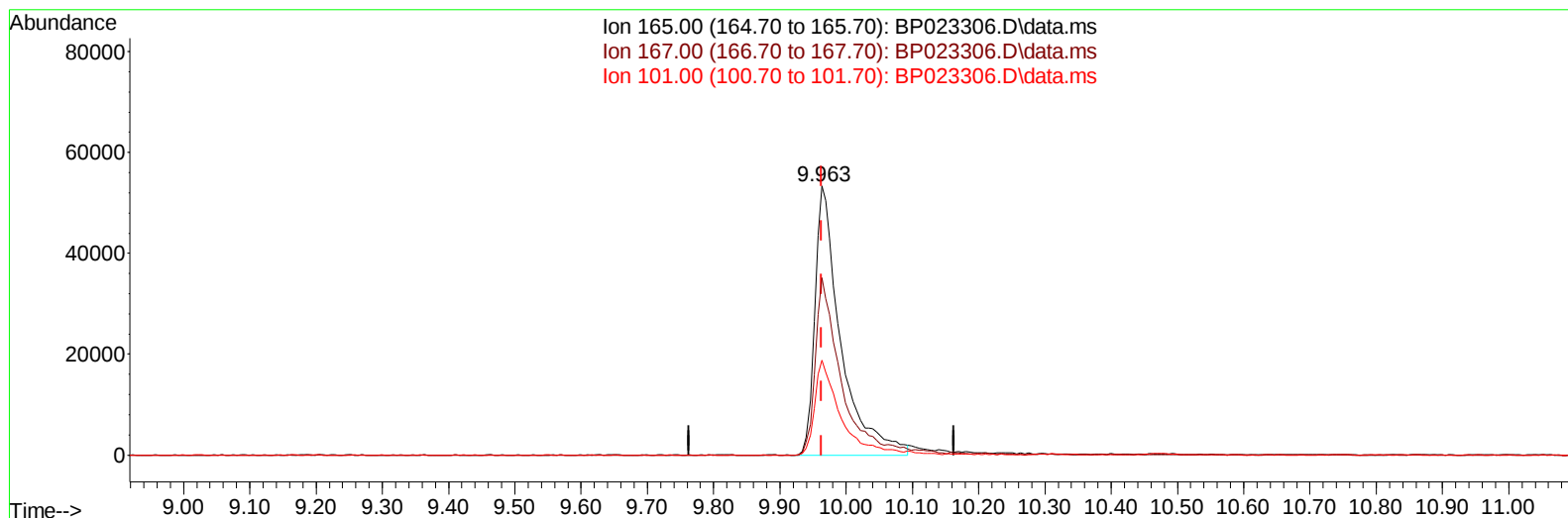
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TIC: BP023306.D\data.ms

(28) 2,4-Dichlorophenol-d3 (S)

9.963min (+ 0.000) 32.51 ng/ul m

response 144763

Ion	Exp%	Act%
165.00	100.00	100.00
167.00	65.80	65.90
101.00	32.60	35.17
0.00	0.00	0.00

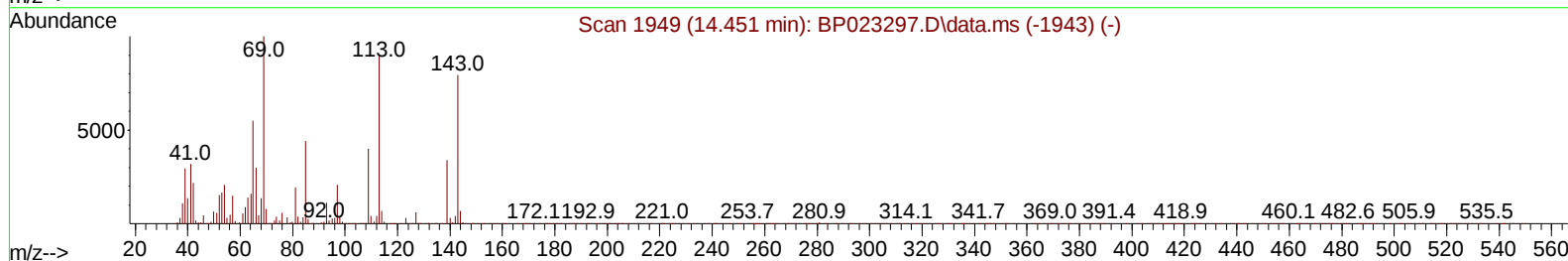
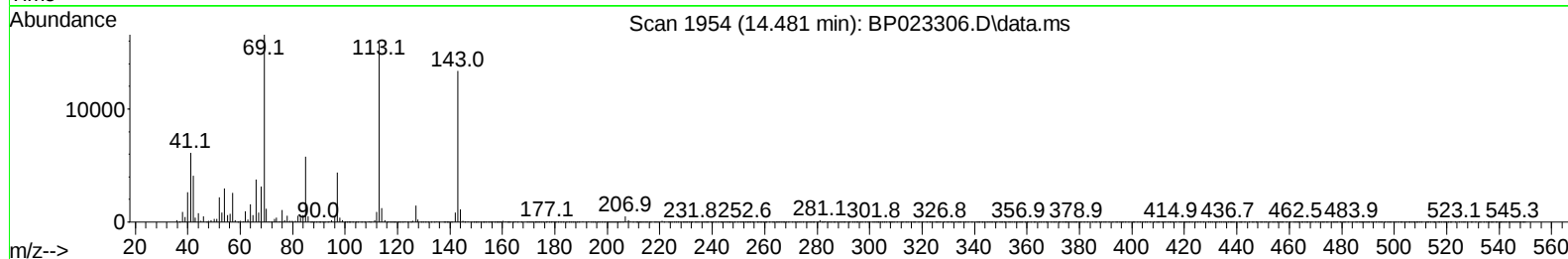
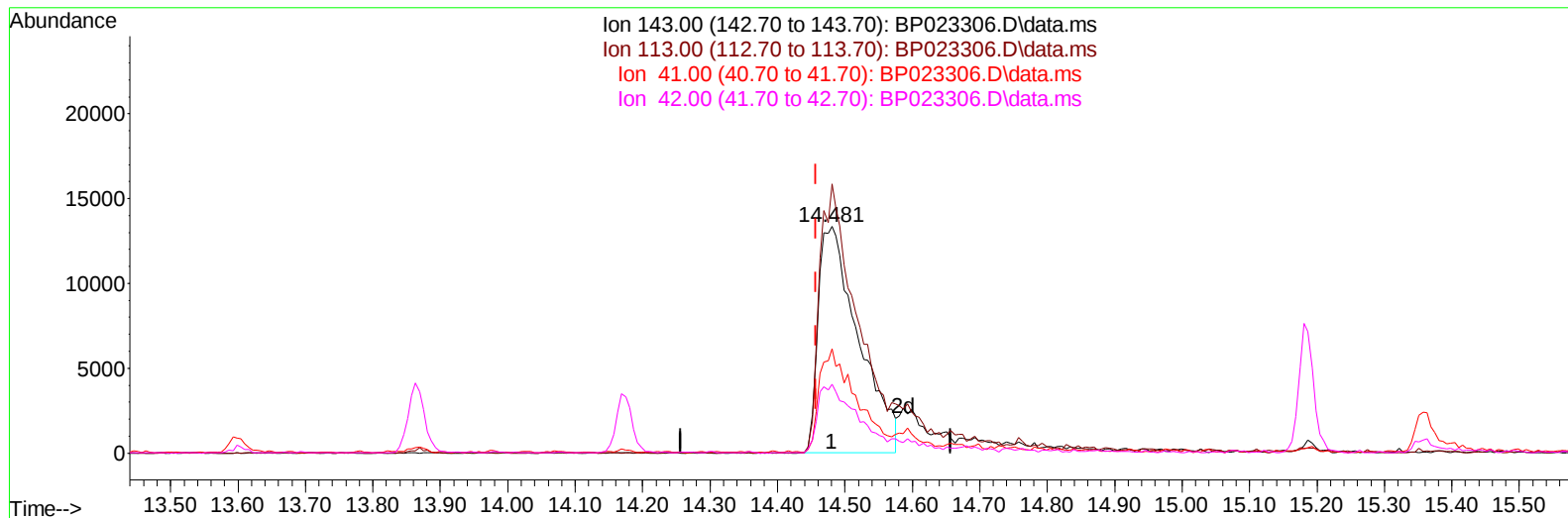
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Instrument :
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TIC: BP023306.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.481min (+ 0.024) 27.92 ng/u1

response 55532

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	107.90	118.69
41.00	39.20	46.01
42.00	25.90	30.52

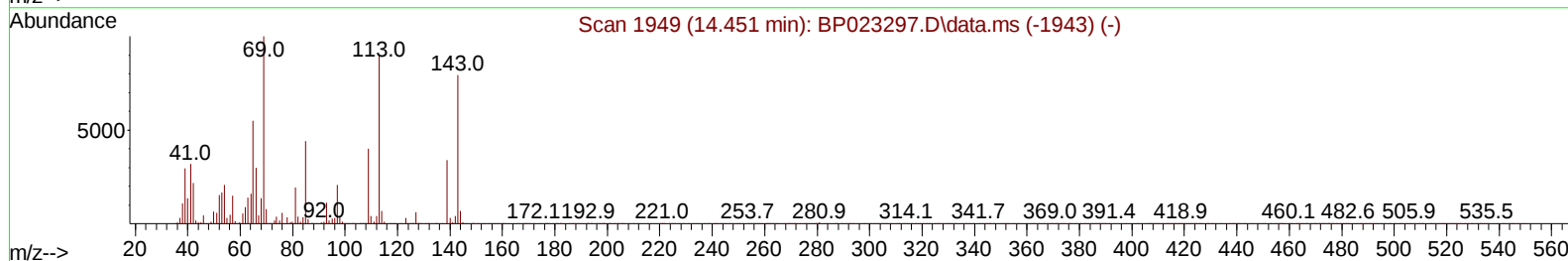
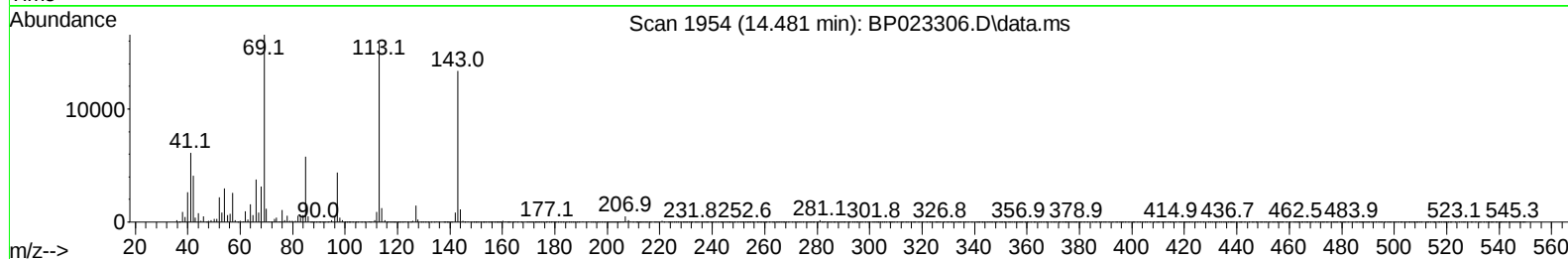
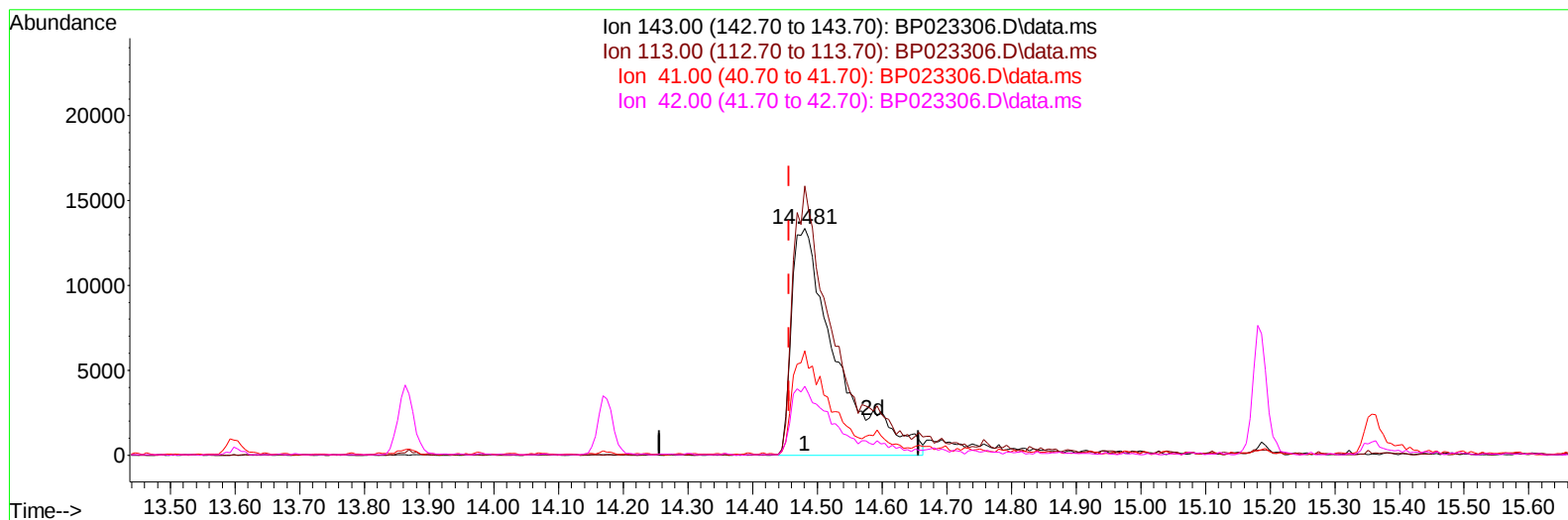
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Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
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Manual IntegrationsAPPROVED

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TIC: BP023306.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.481min (+ 0.024) 32.13 ng/ul m

response 63910

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	107.90	118.69
41.00	39.20	46.01
42.00	25.90	30.52

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Instrument :
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 ClientSampleId :
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Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 11/25/2024
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Quant Time: Nov 22 23:47:36 2024
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.558	152	70832	20.000	ng/ul	-0.01
20) Naphthalene-d8	10.310	136	256169	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.175	164	183647	20.000	ng/ul	-0.01
64) Phenanthrene-d10	16.987	188	447961	20.000	ng/ul	-0.01
79) Chrysene-d12	21.439	240	534394	20.000	ng/ul	0.02
88) Perylene-d12	24.663	264	639498	20.000	ng/ul	0.03
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.134	96	12654	8.587	ng/uL	0.00
4) Pyridine-d5	3.546	84	150114	34.287	ng/ul	0.00
7) Phenol-d5	6.769	99	183000	34.945	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.911	67	120428	39.142	ng/ul	0.00
11) 2-Chlorophenol-d4	7.111	132	139916	36.481	ng/ul	0.00
15) 4-Methylphenol-d8	8.275	113	142280	33.016	ng/ul	0.00
21) Nitrobenzene-d5	8.705	128	66635	37.364	ng/ul	0.00
24) 2-Nitrophenol-d4	9.416	143	76987	36.210	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	9.963	165	144763m	32.513	ng/ul	0.00
31) 4-Chloroaniline-d4	10.469	131	168377	31.699	ng/ul	0.00
46) Dimethylphthalate-d6	13.598	166	515562	38.951	ng/ul	0.01
49) Acenaphthylene-d8	13.863	160	524431	37.973	ng/ul	-0.01
54) 4-Nitrophenol-d4	14.481	143	63910m	32.130	ng/ul	0.02
60) Fluorene-d10	15.187	176	442832	38.160	ng/ul	-0.02
65) 4,6-Dinitro-2-methylph...	15.357	200	76159	28.380	ng/ul	0.00
73) Anthracene-d10	17.087	188	755077	39.810	ng/ul	-0.01
81) Pyrene-d10	19.486	212	988500	40.159	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.439	264	1198857	39.356	ng/ul	0.02

Target Compounds Qvalue

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

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