

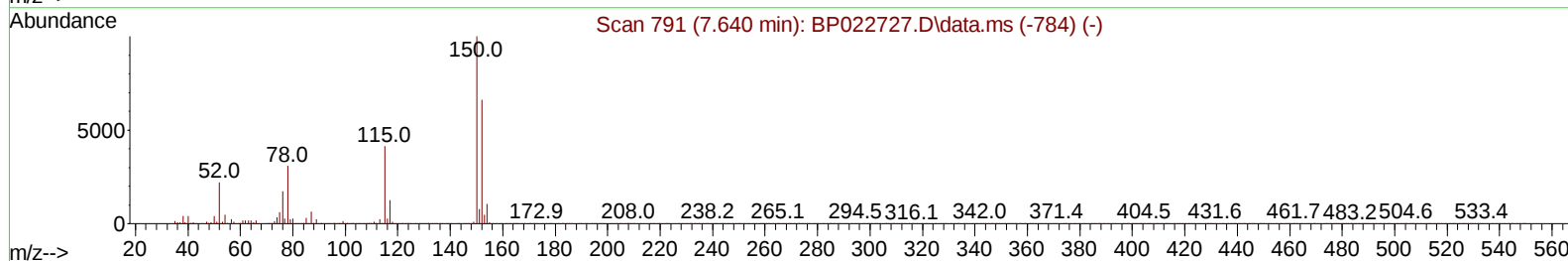
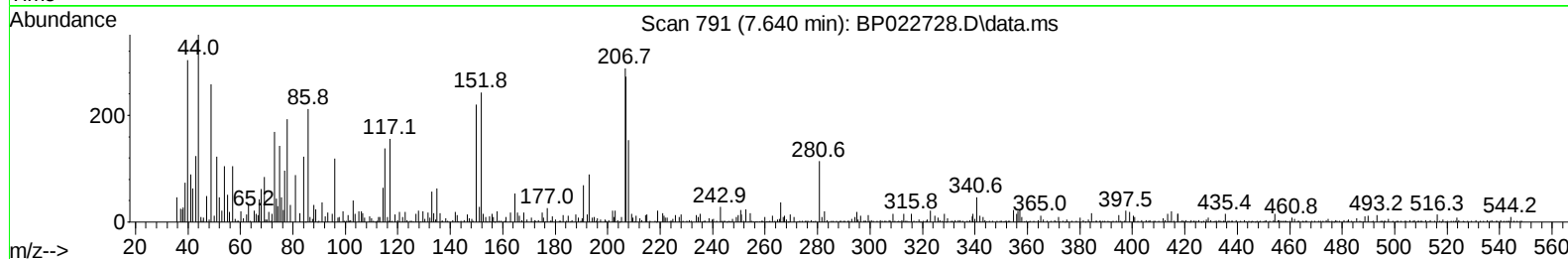
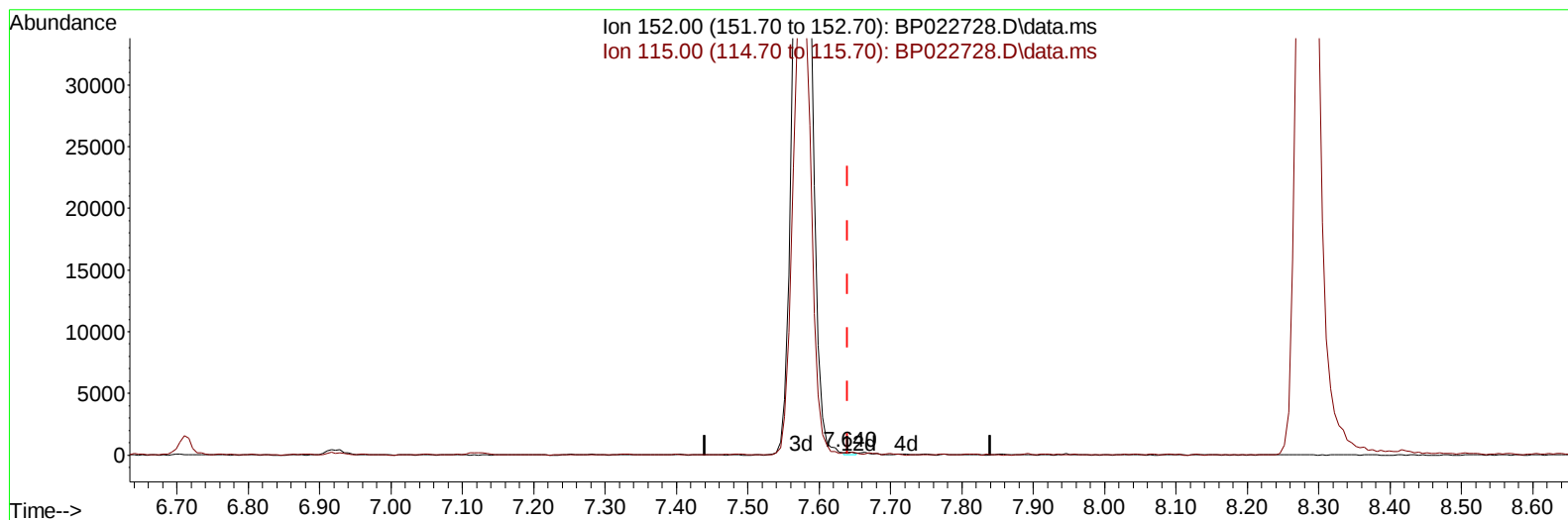
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102924\
 Data File : BP022728.D
 Acq On : 30 Oct 2024 09:04
 Operator : RC/JU
 Sample : PB164416BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK416

Manual IntegrationsAPPROVED

Quant Time: Oct 30 10:57:22 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100724.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Oct 30 10:56:39 2024
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 10/30/2024
 Supervised By :mohammad ahmed 11/04/2024



TIC: BP022728.D\data.ms

(1) 1,4-Dichlorobenzene-d4 (I)

7.640min (+ 0.000) 20.00 ng/u1

response 224

Ion	Exp%	Act%
152.00	100.00	100.00
115.00	59.50	53.28
0.00	0.00	0.00
0.00	0.00	0.00

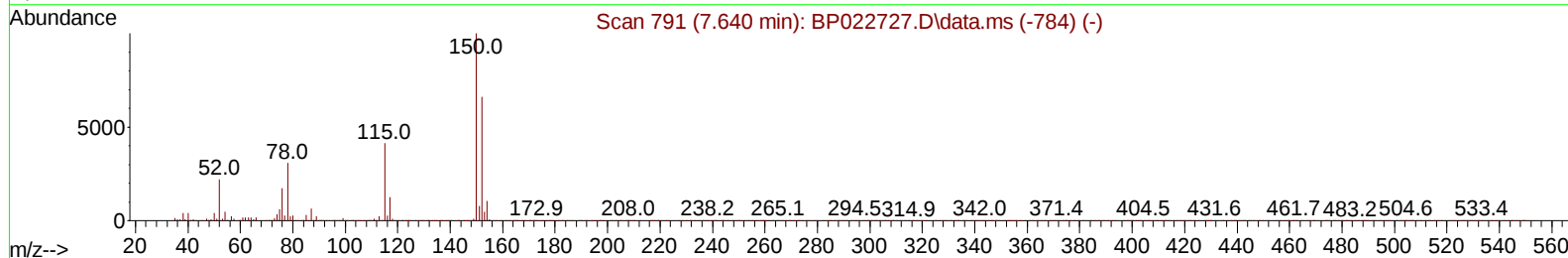
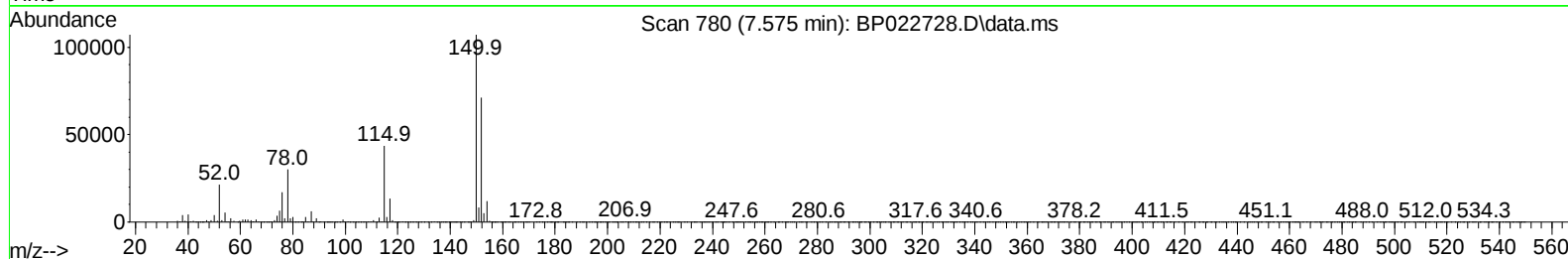
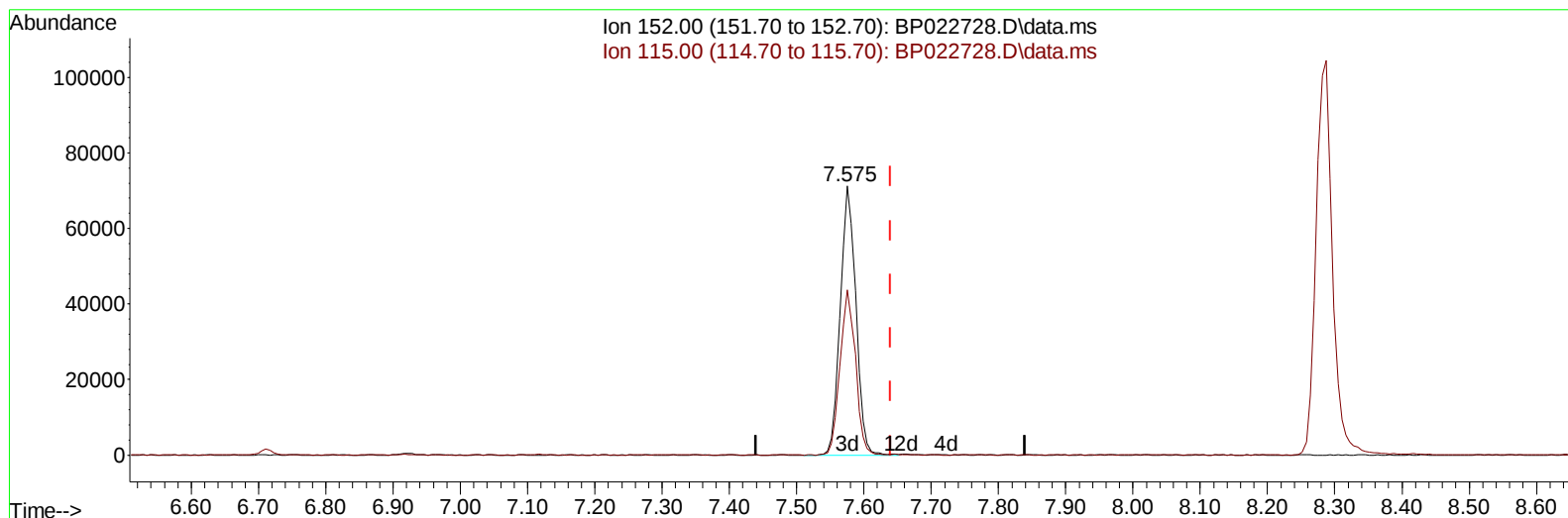
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TIC: BP022728.D\data.ms

(1) 1,4-Dichlorobenzene-d4 (I)

7.575min (-0.065) 20.00 ng/u1 m

response 114797

Ion	Exp%	Act%
152.00	100.00	100.00
115.00	59.50	61.36
0.00	0.00	0.00
0.00	0.00	0.00

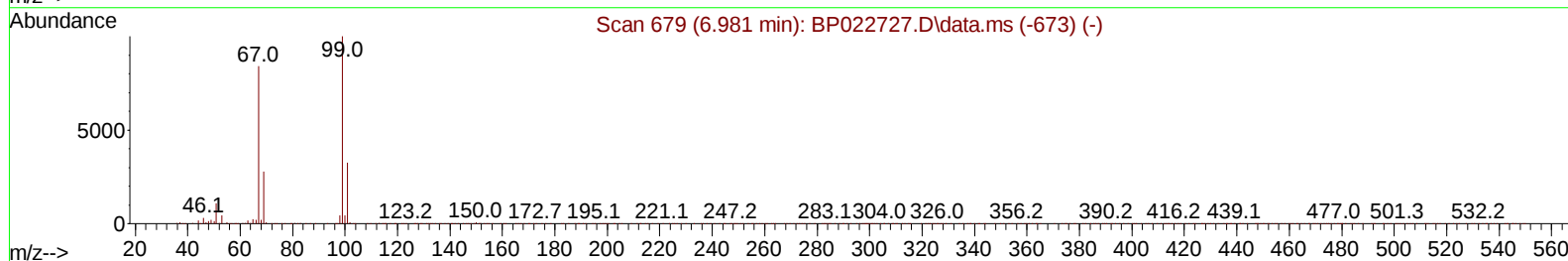
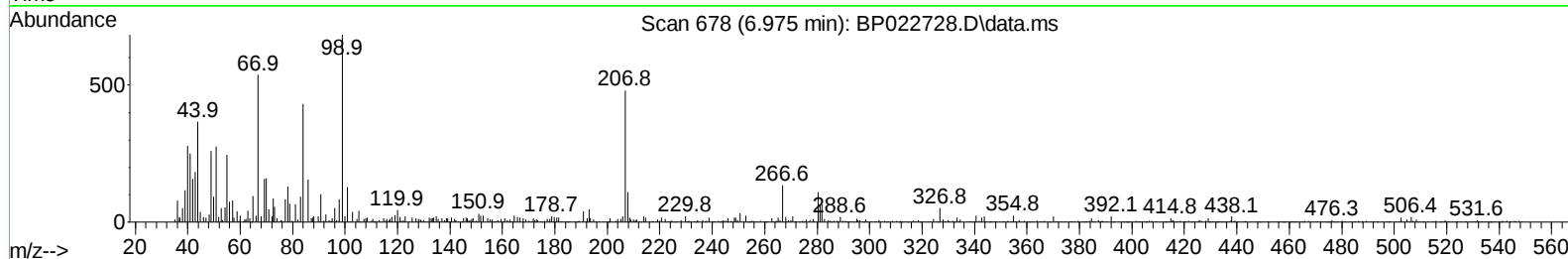
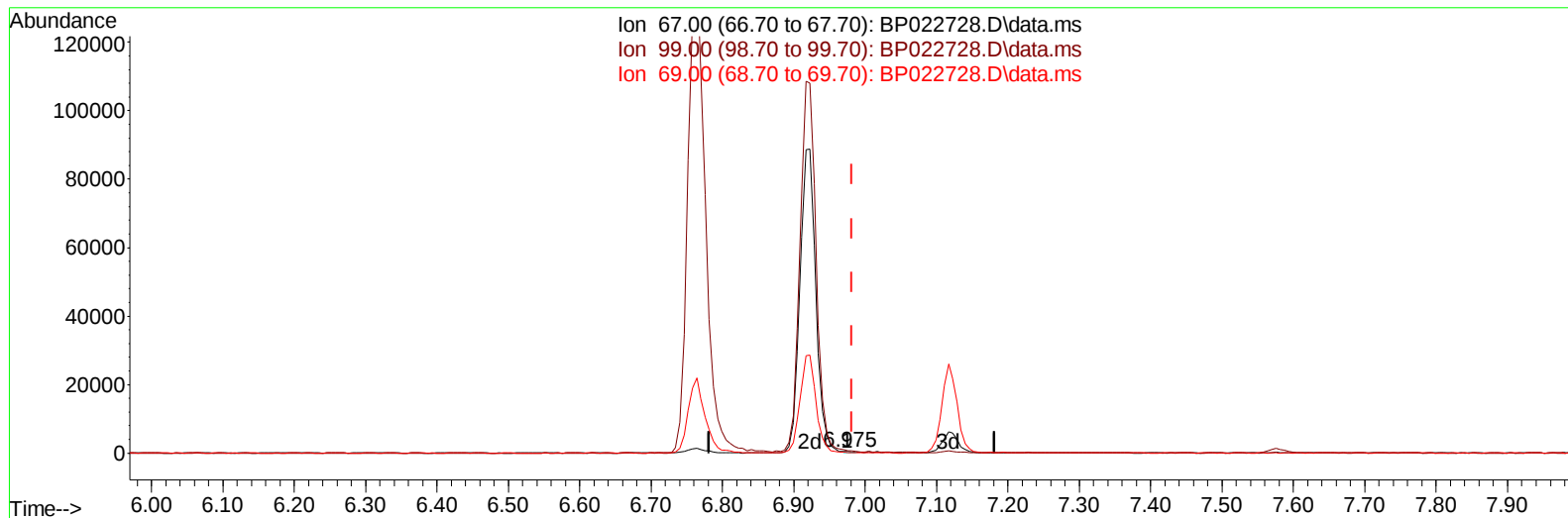
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Instrument :
 BNA_P
 ClientSampleId :
 SBLK416

Manual IntegrationsAPPROVED

Quant Time: Oct 30 11:06:40 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100724.MA.M
 Quant Title : SVOA CALIBRATION
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TIC: BP022728.D\data.ms

(9) Bis-(2-Chloroethyl)ether-d8 (S)

6.975min (-0.006) 0.04 ng/u1

response 211

Ion	Exp%	Act%
67.00	100.00	100.00
99.00	117.00	127.37
69.00	32.90	29.24
0.00	0.00	0.00

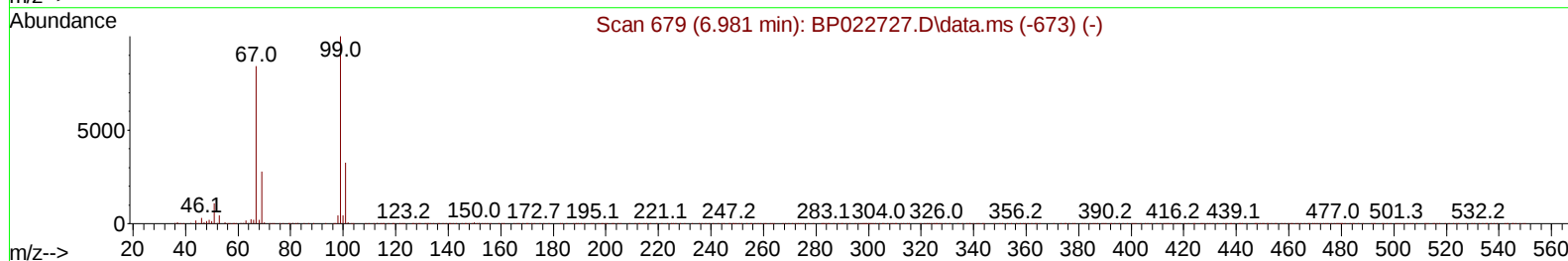
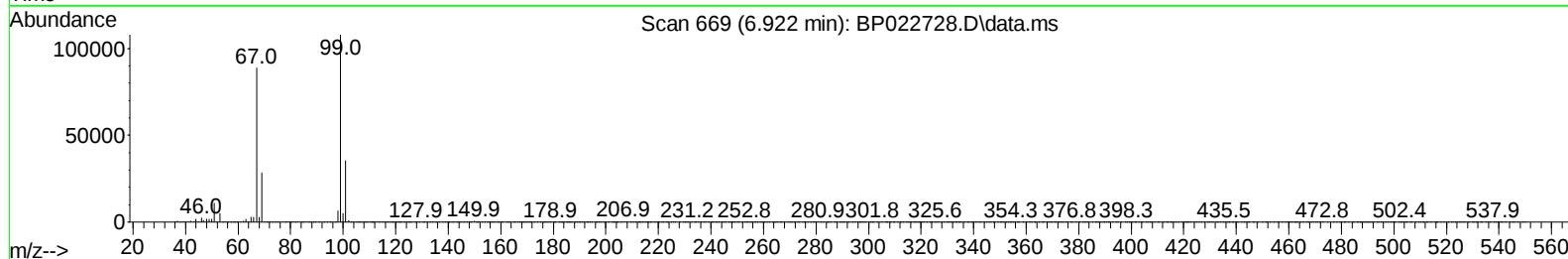
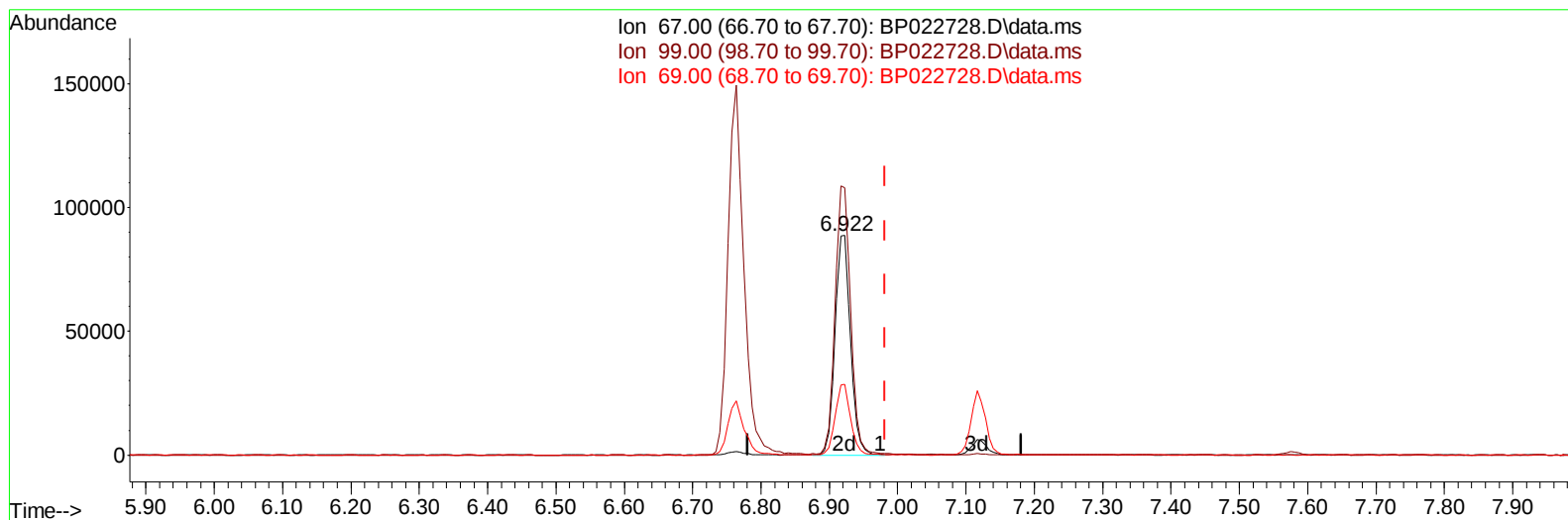
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TIC: BP022728.D\data.ms

(9) Bis-(2-Chloroethyl)ether-d8 (S)

6.922min (-0.059) 24.50 ng/u1 m

response 139159

Ion	Exp%	Act%
67.00	100.00	100.00
99.00	117.00	121.60
69.00	32.90	32.31
0.00	0.00	0.00

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Instrument :
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Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 10/30/2024
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Quant Time: Oct 30 11:07:32 2024
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.575	152	114797m	20.000	ng/ul	-0.06
20) Naphthalene-d8	10.346	136	425311	20.000	ng/ul	-0.05
38) Acenaphthene-d10	14.234	164	343084	20.000	ng/ul	-0.03
64) Phenanthrene-d10	17.063	188	835190	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.498	240	944688	20.000	ng/ul	# 0.00
88) Perylene-d12	24.745	264	1082108	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.146	96	13395	4.534	ng/uL	-0.04
4) Pyridine-d5	3.546	84	197596	24.365	ng/ul	-0.05
7) Phenol-d5	6.764	99	241423	24.984	ng/ul	-0.07
9) Bis-(2-Chloroethyl)eth...	6.922	67	139159m	24.502	ng/ul	-0.06
11) 2-Chlorophenol-d4	7.117	132	185444	27.045	ng/ul	-0.06
15) 4-Methylphenol-d8	8.287	113	197277	25.581	ng/ul	-0.06
21) Nitrobenzene-d5	8.722	128	91394	27.947	ng/ul	-0.06
24) 2-Nitrophenol-d4	9.440	143	109261	28.859	ng/ul	-0.06
28) 2,4-Dichlorophenol-d3	9.981	165	217647	27.048	ng/ul	-0.05
31) 4-Chloroaniline-d4	10.493	131	246901	25.967	ng/ul	-0.05
46) Dimethylphthalate-d6	13.646	166	709043	26.012	ng/ul	-0.02
49) Acenaphthylene-d8	13.922	160	753220	26.016	ng/ul	-0.03
54) 4-Nitrophenol-d4	14.481	143	94954	20.764	ng/ul	-0.03
60) Fluorene-d10	15.257	176	621634	26.292	ng/ul	-0.02
65) 4,6-Dinitro-2-methylph...	15.410	200	109772	19.984	ng/ul	0.00
73) Anthracene-d10	17.169	188	1064635	27.097	ng/ul	0.00
81) Pyrene-d10	19.545	212	1397957	27.983	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.516	264	1663494	28.785	ng/ul	0.00

Target Compounds Qvalue

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

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