

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100423\
 Data File : BP017565.D
 Acq On : 05 Oct 2023 14:48
 Operator : MA/JU
 Sample : 04679-12
 Misc :
 ALS Vial : 39 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 PMW-7S(20230927)

Quant Time: Oct 06 02:40:21 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100423.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Oct 04 16:33:47 2023
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.940	152	114697	20.000	ng/u1	0.00
20) Naphthalene-d8	10.775	136	469410	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.634	164	305581	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.422	188	715709	20.000	ng/u1	0.00
79) Chrysene-d12	21.557	240	735585	20.000	ng/u1	0.00
88) Perylene-d12	24.139	264	837149	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.287	96	11445	3.989	ng/uL	0.00
4) Pyridine-d5	3.728	84	88992	11.851	ng/u1	0.00
7) Phenol-d5	7.099	99	78209	8.104	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.281	67	235314	39.961	ng/u1	0.00
11) 2-Chlorophenol-d4	7.458	132	237516	31.258	ng/u1	0.00
15) 4-Methylphenol-d8	8.664	113	155149	20.083	ng/u1	0.00
21) Nitrobenzene-d5	9.146	128	152697	42.283	ng/u1	0.00
24) 2-Nitrophenol-d4	9.863	143	164920	40.490	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.393	165	294177	38.958	ng/u1	0.00
31) 4-Chloroaniline-d4	10.946	131	363406	33.784	ng/u1	0.00
46) Dimethylphthalate-d6	14.046	166	1058040	44.388	ng/u1	0.00
49) Acenaphthylene-d8	14.328	160	1118326	42.548	ng/u1	0.00
54) 4-Nitrophenol-d4	14.887	143	30049	7.199	ng/u1	0.01
60) Fluorene-d10	15.634	176	911590	45.088	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.775	200	163990	36.753	ng/u1	0.00
73) Anthracene-d10	17.522	188	1562542	47.165	ng/u1	0.00
81) Pyrene-d10	19.775	212	2054536	50.265	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.969	264	2126112	50.325	ng/u1	0.00

Target Compounds	Qvalue

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

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