

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101122\
 Data File : BP012059.D
 Acq On : 12 Oct 2022 02:01
 Operator : CG/JU
 Sample : N4991-14
 Misc :
 ALS Vial : 27 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0BP3

Quant Time: Oct 12 02:41:27 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP101122.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Oct 11 23:54:43 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.857	152	218466	20.000	ng/u1	0.00
20) Naphthalene-d8	10.663	136	880666	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.504	164	499269	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.263	188	1040234	20.000	ng/u1	0.00
79) Chrysene-d12	21.345	240	1150710	20.000	ng/u1	0.00
88) Perylene-d12	23.762	264	1191918	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.281	96	20779	3.750	ng/uL	0.00
4) Pyridine-d5	3.716	84	52705	3.577	ng/u1	0.01
7) Phenol-d5	7.022	99	328041	18.670	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.187	67	210626	21.119	ng/u1	0.00
11) 2-Chlorophenol-d4	7.387	132	290710	20.844	ng/u1	0.00
15) 4-Methylphenol-d8	8.569	113	111636	8.272	ng/u1	0.00
21) Nitrobenzene-d5	9.022	128	134902	21.689	ng/u1	0.00
24) 2-Nitrophenol-d4	9.745	143	140217	22.233	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.286	165	239617	19.604	ng/u1	0.00
31) 4-Chloroaniline-d4	10.810	131	242984	13.490	ng/u1	0.00
46) Dimethylphthalate-d6	13.916	166	758658	23.993	ng/u1	0.00
49) Acenaphthylene-d8	14.198	160	878447	22.467	ng/u1	0.00
54) 4-Nitrophenol-d4	14.710	143	114346	18.420	ng/u1	0.00
60) Fluorene-d10	15.498	176	727104	25.416	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.621	200	65341	11.896	ng/u1	0.00
73) Anthracene-d10	17.357	188	1081216	23.936	ng/u1	0.00
81) Pyrene-d10	19.592	212	1494553	25.352	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.603	264	1480681	25.555	ng/u1	-0.01

Target Compounds Qvalue

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

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