

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101122\
 Data File : BP012052.D
 Acq On : 11 Oct 2022 22:00
 Operator : CG/JU
 Sample : N4991-03
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0BF2

Quant Time: Oct 12 00:24:50 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	238693	20.000	ng/ul	0.00
20) Naphthalene-d8	10.663	136	946854	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.504	164	539677	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.263	188	1125938	20.000	ng/ul	0.00
79) Chrysene-d12	21.351	240	1230234	20.000	ng/ul	0.00
88) Perylene-d12	23.768	264	1362262	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.281	96	26420	4.364	ng/uL	0.00
4) Pyridine-d5	3.722	84	43856	2.724	ng/ul	0.02
7) Phenol-d5	7.022	99	272510	14.195	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.187	67	250509	22.990	ng/ul	0.00
11) 2-Chlorophenol-d4	7.387	132	325718	21.375	ng/ul	0.00
15) 4-Methylphenol-d8	8.569	113	145203	9.848	ng/ul	0.00
21) Nitrobenzene-d5	9.022	128	163063	24.384	ng/ul	0.00
24) 2-Nitrophenol-d4	9.746	143	170688	25.173	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.287	165	292297	22.243	ng/ul	0.00
31) 4-Chloroaniline-d4	10.810	131	344547	17.792	ng/ul	0.00
46) Dimethylphthalate-d6	13.916	166	859932	25.160	ng/ul	0.00
49) Acenaphthylene-d8	14.198	160	1029993	24.370	ng/ul	0.00
54) 4-Nitrophenol-d4	14.716	143	68834	10.258	ng/ul	0.00
60) Fluorene-d10	15.498	176	800607	25.890	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.628	200	75167	12.643	ng/ul	0.00
73) Anthracene-d10	17.363	188	1226306	25.082	ng/ul	0.00
81) Pyrene-d10	19.598	212	1597375	25.345	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.610	264	1735444	26.206	ng/ul	0.00

Target Compounds Qvalue

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

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