

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101122\
 Data File : BP012060.D
 Acq On : 12 Oct 2022 02:35
 Operator : CG/JU
 Sample : N4991-12
 Misc :
 ALS Vial : 28 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0BL5

Quant Time: Oct 12 03:38:48 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	231292	20.000	ng/u1	0.00
20) Naphthalene-d8	10.663	136	933411	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.504	164	515732	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.263	188	1036273	20.000	ng/u1	0.00
79) Chrysene-d12	21.345	240	1130635	20.000	ng/u1	0.00
88) Perylene-d12	23.768	264	1255486	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.275	96	23891	4.073	ng/uL	0.00
4) Pyridine-d5	0.000	84	0d	0.000	ng/u1	
7) Phenol-d5	7.022	99	172663	9.282	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.187	67	259324	24.560	ng/u1	0.00
11) 2-Chlorophenol-d4	7.387	132	265983	18.014	ng/u1	0.00
15) 4-Methylphenol-d8	8.569	113	132914	9.303	ng/u1	0.00
21) Nitrobenzene-d5	9.022	128	167560	25.418	ng/u1	0.00
24) 2-Nitrophenol-d4	9.746	143	173272	25.922	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.287	165	263992	20.378	ng/u1	0.00
31) 4-Chloroaniline-d4	10.804	131	297834	15.601	ng/u1	0.00
46) Dimethylphthalate-d6	13.916	166	948877	29.051	ng/u1	0.00
49) Acenaphthylene-d8	14.198	160	1111619	27.523	ng/u1	0.00
54) 4-Nitrophenol-d4	14.716	143	33150	5.170	ng/u1	0.00
60) Fluorene-d10	15.498	176	867103	29.342	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.622	200	62900	11.495	ng/u1	0.00
73) Anthracene-d10	17.357	188	1307448	29.055	ng/u1	0.00
81) Pyrene-d10	19.592	212	1646033	28.418	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.610	264	1720417	28.189	ng/u1	0.00

Target Compounds Qvalue

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

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