

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP090823\
 Data File : BP017046.D
 Acq On : 09 Sep 2023 11:42
 Operator : MA/JU
 Sample : 04166-13
 Misc :
 ALS Vial : 57 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BBHH3

Quant Time: Sep 10 00:10:08 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP090523.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Sat Sep 09 01:41:21 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.969	152	316937	20.000	ng/u1	0.00
20) Naphthalene-d8	10.799	136	1366021	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.628	164	806886	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.392	188	1729424	20.000	ng/u1	0.00
79) Chrysene-d12	21.492	240	1348858	20.000	ng/u1	0.00
88) Perylene-d12	24.022	264	1361802	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.323	96	48158	5.748	ng/uL	0.00
4) Pyridine-d5	3.764	84	168331	7.031	ng/u1	0.00
7) Phenol-d5	7.117	99	248528	8.583	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.311	67	771236	42.582	ng/u1	0.00
11) 2-Chlorophenol-d4	7.493	132	670839	32.011	ng/u1	0.00
15) 4-Methylphenol-d8	8.675	113	456656	20.701	ng/u1	-0.01
21) Nitrobenzene-d5	9.169	128	431316	40.699	ng/u1	0.00
24) 2-Nitrophenol-d4	9.881	143	443614	38.615	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.405	165	734748	34.643	ng/u1	0.00
31) 4-Chloroaniline-d4	10.958	131	922819	30.158	ng/u1	0.00
46) Dimethylphthalate-d6	14.040	166	2777909	46.250	ng/u1	0.00
49) Acenaphthylene-d8	14.328	160	2994133	43.195	ng/u1	0.00
54) 4-Nitrophenol-d4	14.840	143	78173	6.663	ng/u1	0.00
60) Fluorene-d10	15.622	176	2339287	45.490	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.751	200	354621	34.073	ng/u1	0.00
73) Anthracene-d10	17.498	188	3836011	48.299	ng/u1	0.00
81) Pyrene-d10	19.734	212	4169565	50.750	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.857	264	3380637	48.933	ng/u1	0.00

Target Compounds Qvalue

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

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