

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP090823\
 Data File : BP017008.D
 Acq On : 08 Sep 2023 12:46
 Operator : MA/JU
 Sample : 04123-10
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BH930

Quant Time: Sep 09 01:01:18 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP090523.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Sep 06 01:52:05 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.975	152	421494	20.000	ng/u1	0.00
20) Naphthalene-d8	10.804	136	1851801	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.634	164	1131880	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.398	188	2430541	20.000	ng/u1	-0.01
79) Chrysene-d12	21.498	240	1962623	20.000	ng/u1	-0.01
88) Perylene-d12	24.033	264	1963020	20.000	ng/u1	-0.02

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.323	96	73948	6.636	ng/uL	0.00
4) Pyridine-d5	3.764	84	495634	15.567	ng/u1	0.00
7) Phenol-d5	7.122	99	357542	9.285	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.316	67	942251	39.119	ng/u1	0.00
11) 2-Chlorophenol-d4	7.493	132	863313	30.977	ng/u1	-0.01
15) 4-Methylphenol-d8	8.681	113	615582	20.984	ng/u1	-0.01
21) Nitrobenzene-d5	9.175	128	543334	37.820	ng/u1	0.00
24) 2-Nitrophenol-d4	9.887	143	573126	36.802	ng/u1	-0.01
28) 2,4-Dichlorophenol-d3	10.410	165	952794	33.139	ng/u1	-0.01
31) 4-Chloroaniline-d4	10.963	131	1339715	32.297	ng/u1	0.00
46) Dimethylphthalate-d6	14.045	166	3338821	39.628	ng/u1	0.00
49) Acenaphthylene-d8	14.334	160	3710853	38.164	ng/u1	0.00
54) 4-Nitrophenol-d4	14.845	143	121954	7.410	ng/u1	-0.01
60) Fluorene-d10	15.628	176	2877792	39.893	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.757	200	454268	31.057	ng/u1	0.00
73) Anthracene-d10	17.504	188	4521352	40.506	ng/u1	0.00
81) Pyrene-d10	19.739	212	5129516	42.909	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.868	264	4241259	42.588	ng/u1	-0.01

Target Compounds	Qvalue

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

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