

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP090823\
 Data File : BP017019.D
 Acq On : 08 Sep 2023 19:06
 Operator : MA/JU
 Sample : 04175-04
 Misc :
 ALS Vial : 28 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 EZYN9

Quant Time: Sep 09 00:45:34 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	502208	20.000	ng/u1	0.00
20) Naphthalene-d8	10.804	136	2185680	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.634	164	1346435	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.398	188	2919851	20.000	ng/u1	-0.01
79) Chrysene-d12	21.498	240	2295520	20.000	ng/u1	-0.01
88) Perylene-d12	24.033	264	2191958	20.000	ng/u1	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.328	96	75863	5.714	ng/uL	0.00
4) Pyridine-d5	3.764	84	525246	13.845	ng/u1	0.00
7) Phenol-d5	7.122	99	429734	9.366	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.316	67	1221385	42.558	ng/u1	0.00
11) 2-Chlorophenol-d4	7.493	132	1103032	33.217	ng/u1	-0.01
15) 4-Methylphenol-d8	8.681	113	766024	21.915	ng/u1	-0.01
21) Nitrobenzene-d5	9.175	128	707126	41.702	ng/u1	0.00
24) 2-Nitrophenol-d4	9.887	143	741004	40.313	ng/u1	-0.01
28) 2,4-Dichlorophenol-d3	10.410	165	1207335	35.578	ng/u1	-0.01
31) 4-Chloroaniline-d4	10.963	131	1788892	36.538	ng/u1	0.00
46) Dimethylphthalate-d6	14.045	166	4229816	42.203	ng/u1	0.00
49) Acenaphthylene-d8	14.334	160	4719008	40.799	ng/u1	0.00
54) 4-Nitrophenol-d4	14.845	143	150138	7.669	ng/u1	-0.01
60) Fluorene-d10	15.628	176	3664998	42.710	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.751	200	578758	32.937	ng/u1	-0.01
73) Anthracene-d10	17.504	188	5946096	44.344	ng/u1	0.00
81) Pyrene-d10	19.739	212	6618716	47.337	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.868	264	5150641	46.318	ng/u1	-0.01

Target Compounds Qvalue

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

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