

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP090823\
 Data File : BP017043.D
 Acq On : 09 Sep 2023 09:59
 Operator : MA/JU
 Sample : PB155198BL
 Misc :
 ALS Vial : 54 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK198

Quant Time: Sep 10 00:01:44 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	355831	20.000	ng/u1	0.00
20) Naphthalene-d8	10.799	136	1520912	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.628	164	915055	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.398	188	1946618	20.000	ng/u1	0.00
79) Chrysene-d12	21.492	240	1547655	20.000	ng/u1	0.00
88) Perylene-d12	24.021	264	1575058	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.322	96	76774	8.161	ng/uL	0.00
4) Pyridine-d5	3.758	84	792297	29.476	ng/u1	0.00
7) Phenol-d5	7.122	99	1232596	37.916	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.311	67	812596	39.962	ng/u1	0.00
11) 2-Chlorophenol-d4	7.493	132	926781	39.391	ng/u1	0.00
15) 4-Methylphenol-d8	8.681	113	953304	38.492	ng/u1	0.00
21) Nitrobenzene-d5	9.169	128	459297	38.926	ng/u1	0.00
24) 2-Nitrophenol-d4	9.887	143	488089	38.160	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.410	165	835803	35.394	ng/u1	0.00
31) 4-Chloroaniline-d4	10.957	131	602199	17.676	ng/u1	0.00
46) Dimethylphthalate-d6	14.045	166	2758624	40.500	ng/u1	0.00
49) Acenaphthylene-d8	14.328	160	3017163	38.382	ng/u1	0.00
54) 4-Nitrophenol-d4	14.840	143	453990	34.121	ng/u1	0.00
60) Fluorene-d10	15.628	176	2281479	39.121	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.751	200	355124	30.315	ng/u1	0.00
73) Anthracene-d10	17.498	188	3607466	40.353	ng/u1	0.00
81) Pyrene-d10	19.733	212	4014808	42.590	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.863	264	3312461	41.455	ng/u1	0.00

Target Compounds Qvalue

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

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