

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP090823\
 Data File : BP017042.D
 Acq On : 09 Sep 2023 09:25
 Operator : MA/JU
 Sample : PB155195BL
 Misc :
 ALS Vial : 53 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK195

Quant Time: Sep 09 23:58:08 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	497575	20.000	ng/u1	0.00
20) Naphthalene-d8	10.799	136	2163860	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.628	164	1334425	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.398	188	2885303	20.000	ng/u1	0.00
79) Chrysene-d12	21.498	240	2300124	20.000	ng/u1	0.00
88) Perylene-d12	24.033	264	2285536	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.322	96	78088	5.936	ng/uL	0.00
4) Pyridine-d5	3.758	84	823648	21.913	ng/u1	0.00
7) Phenol-d5	7.122	99	1296454	28.520	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.311	67	859191	30.217	ng/u1	0.00
11) 2-Chlorophenol-d4	7.493	132	983118	29.882	ng/u1	0.00
15) 4-Methylphenol-d8	8.681	113	1014552	29.295	ng/u1	0.00
21) Nitrobenzene-d5	9.169	128	490086	29.194	ng/u1	0.00
24) 2-Nitrophenol-d4	9.887	143	540056	29.677	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.410	165	910157	27.091	ng/u1	0.00
31) 4-Chloroaniline-d4	10.957	131	652284	13.457	ng/u1	0.00
46) Dimethylphthalate-d6	14.045	166	3028225	30.486	ng/u1	0.00
49) Acenaphthylene-d8	14.328	160	3313909	28.909	ng/u1	0.00
54) 4-Nitrophenol-d4	14.839	143	511495	26.362	ng/u1	0.00
60) Fluorene-d10	15.622	176	2523855	29.676	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.751	200	402253	23.166	ng/u1	0.00
73) Anthracene-d10	17.498	188	4010937	30.270	ng/u1	0.00
81) Pyrene-d10	19.733	212	4432101	31.635	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.863	264	3552114	30.635	ng/u1	0.00

Target Compounds Qvalue

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

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