

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP091823\
 Data File : BP017211.D
 Acq On : 18 Sep 2023 18:33
 Operator : MA/JU
 Sample : PB155579BL
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK579

Quant Time: Sep 18 22:34:08 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP091223.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Sat Sep 16 03:18:48 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.946	152	257155	20.000	ng/ul	0.00
20) Naphthalene-d8	10.769	136	1091361	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.604	164	681819	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.375	188	1444773	20.000	ng/ul	0.00
79) Chrysene-d12	21.475	240	982337	20.000	ng/ul	0.00
88) Perylene-d12	23.998	264	988555	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.305	96	33236	5.431	ng/uL	0.00
4) Pyridine-d5	3.740	84	501442	28.645	ng/ul	0.00
7) Phenol-d5	7.093	99	601968	28.788	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.287	67	383179	29.319	ng/ul	0.00
11) 2-Chlorophenol-d4	7.464	132	474521	28.447	ng/ul	0.00
15) 4-Methylphenol-d8	8.652	113	474732	27.394	ng/ul	0.00
21) Nitrobenzene-d5	9.140	128	230199	30.123	ng/ul	0.00
24) 2-Nitrophenol-d4	9.858	143	253643	31.090	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.381	165	448636	28.434	ng/ul	0.00
31) 4-Chloroaniline-d4	10.934	131	673038	28.301	ng/ul	0.00
46) Dimethylphthalate-d6	14.016	166	1477817	30.458	ng/ul	0.00
49) Acenaphthylene-d8	14.304	160	1601858	29.096	ng/ul	0.00
54) 4-Nitrophenol-d4	14.822	143	193868	20.674	ng/ul	0.00
60) Fluorene-d10	15.598	176	1253875	29.888	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.728	200	168572	20.884	ng/ul	0.00
73) Anthracene-d10	17.475	188	1951481	31.791	ng/ul	0.00
81) Pyrene-d10	19.716	212	2075480	40.286	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.833	264	1486575	31.861	ng/ul	0.00

Target Compounds	Qvalue

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

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