

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092723\
 Data File : BP017391.D
 Acq On : 27 Sep 2023 11:29
 Operator : MA/JU
 Sample : PB155856BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK856

Quant Time: Sep 27 22:37:51 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092023.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Sep 26 04:52:29 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.916	152	115847	20.000	ng/u1	-0.01
20) Naphthalene-d8	10.740	136	530964	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.581	164	339440	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.351	188	753377	20.000	ng/u1	0.00
79) Chrysene-d12	21.457	240	665638	20.000	ng/u1	0.00
88) Perylene-d12	23.968	264	704677	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.293	96	22938	8.133	ng/uL	0.00
4) Pyridine-d5	3.722	84	326635	41.426	ng/u1	0.00
7) Phenol-d5	7.075	99	392556	41.259	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.263	67	245392	42.737	ng/u1	0.00
11) 2-Chlorophenol-d4	7.440	132	316759	41.976	ng/u1	0.00
15) 4-Methylphenol-d8	8.628	113	310069	41.918	ng/u1	-0.01
21) Nitrobenzene-d5	9.116	128	151511	39.471	ng/u1	-0.01
24) 2-Nitrophenol-d4	9.828	143	162468	38.980	ng/u1	-0.01
28) 2,4-Dichlorophenol-d3	10.357	165	303246	38.510	ng/u1	0.00
31) 4-Chloroaniline-d4	10.904	131	453251	39.735	ng/u1	-0.01
46) Dimethylphthalate-d6	13.998	166	991085	40.758	ng/u1	0.00
49) Acenaphthylene-d8	14.281	160	1079345	38.594	ng/u1	0.00
54) 4-Nitrophenol-d4	14.816	143	105383	24.876	ng/u1	0.00
60) Fluorene-d10	15.581	176	856965	40.936	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.710	200	84603	19.924	ng/u1	0.00
73) Anthracene-d10	17.451	188	1374549	41.138	ng/u1	0.00
81) Pyrene-d10	19.698	212	1582062	43.951	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.804	264	1432568	42.224	ng/u1	0.00

Target Compounds Qvalue

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

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