

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092823\
 Data File : BP017420.D
 Acq On : 28 Sep 2023 19:12
 Operator : MA/JU
 Sample : PB155751BL
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK751

Quant Time: Sep 28 23:30:47 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092023.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Sep 28 23:22:11 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.963	152	147576	20.000	ng/u1	0.00
20) Naphthalene-d8	10.804	136	644354	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.663	164	417911	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.451	188	974870	20.000	ng/u1	0.00
79) Chrysene-d12	21.580	240	885103	20.000	ng/u1	0.00
88) Perylene-d12	24.168	264	1059694	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.299	96	24786	6.899	ng/uL	0.00
4) Pyridine-d5	3.740	84	344696	34.317	ng/u1	0.00
7) Phenol-d5	7.116	99	431525	35.603	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.305	67	250303	34.220	ng/u1	0.00
11) 2-Chlorophenol-d4	7.481	132	350979	36.511	ng/u1	0.00
15) 4-Methylphenol-d8	8.687	113	339138	35.990	ng/u1	0.00
21) Nitrobenzene-d5	9.175	128	172815	37.098	ng/u1	0.00
24) 2-Nitrophenol-d4	9.893	143	190300	37.623	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.422	165	330710	34.607	ng/u1	0.00
31) 4-Chloroaniline-d4	10.975	131	478163	34.542	ng/u1	0.00
46) Dimethylphthalate-d6	14.075	166	1047140	34.977	ng/u1	0.00
49) Acenaphthylene-d8	14.357	160	1174322	34.105	ng/u1	0.00
54) 4-Nitrophenol-d4	14.898	143	168421	32.291	ng/u1	0.00
60) Fluorene-d10	15.663	176	905905	35.148	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.810	200	112397	20.456	ng/u1	0.00
73) Anthracene-d10	17.551	188	1551193	35.877	ng/u1	0.00
81) Pyrene-d10	19.804	212	1791220	37.423	ng/u1	0.00
92) Benzo(a)pyrene-d12	24.004	264	1816936	35.611	ng/u1	0.00

Target Compounds Qvalue

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

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