

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092923\
 Data File : BP017467.D
 Acq On : 30 Sep 2023 04:06
 Operator : MA/JU
 Sample : PB155871BL
 Misc :
 ALS Vial : 33 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK871

Quant Time: Sep 30 04:40:40 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092023.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Sat Sep 30 04:02:20 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.952	152	197172	20.000	ng/ul	0.00
20) Naphthalene-d8	10.787	136	791558	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.639	164	467020	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.428	188	997750	20.000	ng/ul	0.00
79) Chrysene-d12	21.557	240	979146	20.000	ng/ul	0.00
88) Perylene-d12	24.127	264	1124774	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.293	96	22165	4.617	ng/uL	0.00
4) Pyridine-d5	3.728	84	317416	23.652	ng/ul	0.00
7) Phenol-d5	7.105	99	392476	24.236	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.293	67	239538	24.511	ng/ul	0.00
11) 2-Chlorophenol-d4	7.469	132	321233	25.011	ng/ul	0.00
15) 4-Methylphenol-d8	8.669	113	315482	25.058	ng/ul	0.00
21) Nitrobenzene-d5	9.157	128	155850	27.235	ng/ul	0.00
24) 2-Nitrophenol-d4	9.875	143	172770	27.805	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.404	165	303563	25.859	ng/ul	0.00
31) 4-Chloroaniline-d4	10.957	131	436605	25.675	ng/ul	0.00
46) Dimethylphthalate-d6	14.057	166	916585	27.397	ng/ul	0.00
49) Acenaphthylene-d8	14.339	160	1035289	26.906	ng/ul	0.00
54) 4-Nitrophenol-d4	14.881	143	119890	20.569	ng/ul	0.00
60) Fluorene-d10	15.645	176	773383	26.851	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.786	200	103257	18.362	ng/ul	0.00
73) Anthracene-d10	17.528	188	1243121	28.093	ng/ul	0.00
81) Pyrene-d10	19.780	212	1482940	28.007	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.957	264	1495958	27.624	ng/ul	0.00

Target Compounds Qvalue

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

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