

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP091523\
 Data File : BP017190.D
 Acq On : 15 Sep 2023 16:27
 Operator : MA/JU
 Sample : PB155306BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK306

Quant Time: Sep 16 03:20:18 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	135656	20.000	ng/ul	0.00
20) Naphthalene-d8	10.775	136	530973	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.604	164	292614	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.375	188	624610	20.000	ng/ul	0.00
79) Chrysene-d12	21.480	240	567059	20.000	ng/ul	0.00
88) Perylene-d12	24.004	264	644839	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.311	96	25202	7.807	ng/uL	0.00
4) Pyridine-d5	3.746	84	226959	24.577	ng/ul	0.00
7) Phenol-d5	7.099	99	380579	34.502	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.287	67	250972	36.402	ng/ul	0.00
11) 2-Chlorophenol-d4	7.469	132	299965	34.088	ng/ul	0.00
15) 4-Methylphenol-d8	8.657	113	282686	30.922	ng/ul	0.00
21) Nitrobenzene-d5	9.146	128	129048	34.709	ng/ul	0.00
24) 2-Nitrophenol-d4	9.863	143	135418	34.116	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.387	165	258298	33.648	ng/ul	0.00
31) 4-Chloroaniline-d4	10.934	131	254464	21.993	ng/ul	0.00
46) Dimethylphthalate-d6	14.022	166	738492	35.465	ng/ul	0.00
49) Acenaphthylene-d8	14.304	160	811894	34.362	ng/ul	0.00
54) 4-Nitrophenol-d4	14.828	143	95420	23.710	ng/ul	0.00
60) Fluorene-d10	15.604	176	659805	36.646	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.734	200	77480	22.203	ng/ul	0.00
73) Anthracene-d10	17.475	188	1045320	39.389	ng/ul	0.00
81) Pyrene-d10	19.716	212	1213467	40.803	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.833	264	1243556	40.859	ng/ul	0.00

Target Compounds Qvalue

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

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