

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP091223\
 Data File : BP017121.D
 Acq On : 12 Sep 2023 17:49
 Operator : MA/JU
 Sample : PB155306BL
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK306

Quant Time: Sep 13 00:56:42 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP091223.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Sep 12 23:27:27 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.952	152	208651	20.000	ng/u1	0.00
20) Naphthalene-d8	10.781	136	823530	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.616	164	475820	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.387	188	937320	20.000	ng/u1	0.00
79) Chrysene-d12	21.486	240	710020	20.000	ng/u1	0.00
88) Perylene-d12	24.022	264	816680	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.311	96	32829	6.612	ng/uL	0.00
4) Pyridine-d5	3.746	84	428360	30.158	ng/u1	0.00
7) Phenol-d5	7.099	99	510206	30.072	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.287	67	322473	30.410	ng/u1	-0.01
11) 2-Chlorophenol-d4	7.469	132	411478	30.402	ng/u1	0.00
15) 4-Methylphenol-d8	8.658	113	392117	27.887	ng/u1	-0.01
21) Nitrobenzene-d5	9.152	128	187555	32.524	ng/u1	0.00
24) 2-Nitrophenol-d4	9.863	143	202252	32.853	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.387	165	366329	30.769	ng/u1	0.00
31) 4-Chloroaniline-d4	10.940	131	529508	29.507	ng/u1	0.00
46) Dimethylphthalate-d6	14.028	166	1076148	31.782	ng/u1	0.00
49) Acenaphthylene-d8	14.316	160	1215778	31.644	ng/u1	0.00
54) 4-Nitrophenol-d4	14.834	143	145052	22.165	ng/u1	0.00
60) Fluorene-d10	15.610	176	937438	32.019	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.740	200	115973	22.146	ng/u1	0.00
73) Anthracene-d10	17.487	188	1349616	33.889	ng/u1	0.00
81) Pyrene-d10	19.728	212	1423822	38.237	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.851	264	1313842	34.085	ng/u1	0.00

Target Compounds Qvalue

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

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