

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110123\
 Data File : BP017903.D
 Acq On : 03 Nov 2023 01:06
 Operator : MA/JU
 Sample : PB156797BL
 Misc :
 ALS Vial : 59 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK797

Quant Time: Nov 03 01:46:49 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP101823.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Oct 19 01:32:15 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.875	152	167312	20.000	ng/u1	-0.01
20) Naphthalene-d8	10.705	136	627251	20.000	ng/u1	-0.02
38) Acenaphthene-d10	14.575	164	355407	20.000	ng/u1	-0.01
64) Phenanthrene-d10	17.369	188	714437	20.000	ng/u1	0.00
79) Chrysene-d12	21.516	240	571414	20.000	ng/u1	0.00
88) Perylene-d12	24.074	264	583592	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.276	96	21196	4.713	ng/uL	0.02
4) Pyridine-d5	3.705	84	270068	22.935	ng/u1	0.01
7) Phenol-d5	7.040	99	378819	27.208	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.222	67	234683	27.294	ng/u1	-0.01
11) 2-Chlorophenol-d4	7.399	132	313501	27.607	ng/u1	-0.01
15) 4-Methylphenol-d8	8.599	113	294981	26.979	ng/u1	-0.01
21) Nitrobenzene-d5	9.087	128	150689	29.648	ng/u1	-0.01
24) 2-Nitrophenol-d4	9.799	143	167120	29.759	ng/u1	-0.01
28) 2,4-Dichlorophenol-d3	10.328	165	279513	25.373	ng/u1	-0.01
31) 4-Chloroaniline-d4	10.881	131	384372	26.661	ng/u1	-0.01
46) Dimethylphthalate-d6	13.993	166	821315	27.528	ng/u1	-0.01
49) Acenaphthylene-d8	14.275	160	921452	27.694	ng/u1	0.00
54) 4-Nitrophenol-d4	14.834	143	97860	23.020	ng/u1	0.00
60) Fluorene-d10	15.581	176	666269	25.834	ng/u1	-0.01
65) 4,6-Dinitro-2-methylph...	15.734	200	64025	13.605	ng/u1	0.00
73) Anthracene-d10	17.469	188	1012534	27.733	ng/u1	0.00
81) Pyrene-d10	19.728	212	1094567	30.553	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.904	264	900007	27.689	ng/u1	0.00

Target Compounds Qvalue

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

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