

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100223\
 Data File : BP017494.D
 Acq On : 03 Oct 2023 01:22
 Operator : MA/JU
 Sample : PB155906BL
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK906

Quant Time: Oct 03 02:07:46 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	7.940	152	140351	20.000	ng/u1	-0.01
20) Naphthalene-d8	10.781	136	590465	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.634	164	339113	20.000	ng/u1	-0.01
64) Phenanthrene-d10	17.422	188	663870	20.000	ng/u1	0.00
79) Chrysene-d12	21.551	240	636380	20.000	ng/u1	0.00
88) Perylene-d12	24.121	264	703286	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.287	96	25018	7.322	ng/uL	0.00
4) Pyridine-d5	3.723	84	370811	38.818	ng/u1	0.00
7) Phenol-d5	7.093	99	466143	40.439	ng/u1	-0.01
9) Bis-(2-Chloroethyl)eth...	7.281	67	288258	41.437	ng/u1	-0.01
11) 2-Chlorophenol-d4	7.458	132	372883	40.786	ng/u1	-0.01
15) 4-Methylphenol-d8	8.658	113	373266	41.651	ng/u1	-0.01
21) Nitrobenzene-d5	9.152	128	187956	44.031	ng/u1	0.00
24) 2-Nitrophenol-d4	9.863	143	203475	43.899	ng/u1	-0.01
28) 2,4-Dichlorophenol-d3	10.393	165	364805	41.659	ng/u1	-0.01
31) 4-Chloroaniline-d4	10.952	131	510598	40.252	ng/u1	0.00
46) Dimethylphthalate-d6	14.051	166	1088663	44.814	ng/u1	0.00
49) Acenaphthylene-d8	14.334	160	1189762	42.583	ng/u1	0.00
54) 4-Nitrophenol-d4	14.869	143	115834	27.369	ng/u1	0.00
60) Fluorene-d10	15.640	176	881379	42.143	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.781	200	115556	30.883	ng/u1	0.00
73) Anthracene-d10	17.522	188	1336574	45.395	ng/u1	0.00
81) Pyrene-d10	19.775	212	1585040	46.058	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.957	264	1531145	45.218	ng/u1	0.00

Target Compounds Qvalue

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

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