

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP091223\
 Data File : BP017161.D
 Acq On : 13 Sep 2023 22:02
 Operator : MA/JU
 Sample : PB155240BL
 Misc :
 ALS Vial : 49 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK240

Quant Time: Sep 14 00:24:32 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	239511	20.000	ng/u1	0.00
20) Naphthalene-d8	10.775	136	1017933	20.000	ng/u1	-0.01
38) Acenaphthene-d10	14.610	164	650449	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.381	188	1495543	20.000	ng/u1	0.00
79) Chrysene-d12	21.480	240	1304388	20.000	ng/u1	0.00
88) Perylene-d12	24.004	264	1255354	20.000	ng/u1	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.311	96	36590	6.420	ng/uL	0.00
4) Pyridine-d5	3.746	84	335847	20.599	ng/u1	0.00
7) Phenol-d5	7.099	99	584622	30.019	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.293	67	380607	31.268	ng/u1	0.00
11) 2-Chlorophenol-d4	7.469	132	463649	29.842	ng/u1	0.00
15) 4-Methylphenol-d8	8.658	113	463309	28.705	ng/u1	-0.01
21) Nitrobenzene-d5	9.152	128	223646	31.376	ng/u1	0.00
24) 2-Nitrophenol-d4	9.863	143	244224	32.094	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.387	165	438659	29.807	ng/u1	0.00
31) 4-Chloroaniline-d4	10.940	131	446817	20.144	ng/u1	0.00
46) Dimethylphthalate-d6	14.028	166	1532896	33.117	ng/u1	0.00
49) Acenaphthylene-d8	14.310	160	1641692	31.258	ng/u1	0.00
54) 4-Nitrophenol-d4	14.828	143	228840	25.580	ng/u1	0.00
60) Fluorene-d10	15.604	176	1307454	32.668	ng/u1	-0.01
65) 4,6-Dinitro-2-methylph...	15.734	200	203691	24.378	ng/u1	-0.01
73) Anthracene-d10	17.481	188	2144942	33.756	ng/u1	0.00
81) Pyrene-d10	19.716	212	2523230	36.885	ng/u1	-0.01
92) Benzo(a)pyrene-d12	23.839	264	2055993	34.700	ng/u1	-0.01

Target Compounds Qvalue

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

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