

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP090823\
 Data File : BP017006.D
 Acq On : 08 Sep 2023 11:37
 Operator : MA/JU
 Sample : PB155065BL
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK065

Quant Time: Sep 09 00:51:45 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP090523.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Sep 06 01:52:05 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.975	152	391333	20.000	ng/u1	0.00
20) Naphthalene-d8	10.805	136	1621685	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.634	164	944658	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.404	188	1890082	20.000	ng/u1	0.00
79) Chrysene-d12	21.498	240	1304549	20.000	ng/u1	-0.01
88) Perylene-d12	24.039	264	1325577	20.000	ng/u1	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.328	96	91054	8.801	ng/uL	0.00
4) Pyridine-d5	3.764	84	1045958	35.383	ng/u1	0.00
7) Phenol-d5	7.122	99	1211227	33.878	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.316	67	793928	35.502	ng/u1	0.00
11) 2-Chlorophenol-d4	7.493	132	917104	35.443	ng/u1	-0.01
15) 4-Methylphenol-d8	8.681	113	917842	33.698	ng/u1	-0.01
21) Nitrobenzene-d5	9.175	128	444119	35.300	ng/u1	0.00
24) 2-Nitrophenol-d4	9.887	143	482976	35.414	ng/u1	-0.01
28) 2,4-Dichlorophenol-d3	10.410	165	826555	32.828	ng/u1	-0.01
31) 4-Chloroaniline-d4	10.963	131	1227780	33.799	ng/u1	0.00
46) Dimethylphthalate-d6	14.046	166	2526793	35.934	ng/u1	0.00
49) Acenaphthylene-d8	14.334	160	2871667	35.387	ng/u1	0.00
54) 4-Nitrophenol-d4	14.845	143	415426	30.244	ng/u1	-0.01
60) Fluorene-d10	15.628	176	2146434	35.652	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.757	200	303121	26.649	ng/u1	0.00
73) Anthracene-d10	17.504	188	3176708	36.598	ng/u1	0.00
81) Pyrene-d10	19.739	212	3307596	41.626	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.869	264	2656128	39.497	ng/u1	-0.01

Target Compounds Qvalue

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

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