

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101923\
 Data File : BP017726.D
 Acq On : 19 Oct 2023 12:24
 Operator : MA/JU
 Sample : PB156344TB 10X
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SLEB344

Quant Time: Oct 20 00:44:36 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	7.887	152	124963	20.000	ng/u1	0.00
20) Naphthalene-d8	10.722	136	468073	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.587	164	286374	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.381	188	634284	20.000	ng/u1	0.00
79) Chrysene-d12	21.522	240	584456	20.000	ng/u1	0.00
88) Perylene-d12	24.074	264	646467	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.258	96	20627	6.141	ng/uL	0.00
4) Pyridine-d5	3.693	84	249896	28.414	ng/u1	0.00
7) Phenol-d5	7.046	99	326479	31.396	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.234	67	216001	33.634	ng/u1	0.00
11) 2-Chlorophenol-d4	7.411	132	289477	34.131	ng/u1	0.00
15) 4-Methylphenol-d8	8.611	113	256771	31.443	ng/u1	0.00
21) Nitrobenzene-d5	9.099	128	129788	34.219	ng/u1	0.00
24) 2-Nitrophenol-d4	9.816	143	147939	35.302	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.340	165	260401	31.677	ng/u1	0.00
31) 4-Chloroaniline-d4	10.893	131	324344	30.149	ng/u1	0.00
46) Dimethylphthalate-d6	14.004	166	866324	36.036	ng/u1	0.00
49) Acenaphthylene-d8	14.287	160	927854	34.609	ng/u1	0.00
54) 4-Nitrophenol-d4	14.834	143	80215	23.418	ng/u1	0.00
60) Fluorene-d10	15.592	176	745355	35.867	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.740	200	115687	27.690	ng/u1	0.00
73) Anthracene-d10	17.481	188	1154632	35.622	ng/u1	0.00
81) Pyrene-d10	19.739	212	1371086	37.418	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.904	264	1319551	36.648	ng/u1	0.00

Target Compounds Qvalue

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

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