

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM040124\
 Data File : BM044858.D
 Acq On : 01 Apr 2024 10:43
 Operator : MA/JU
 Sample : PB159895BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK895

Quant Time: Apr 01 23:15:19 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM030524.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Sun Mar 31 13:54:26 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.663	152	108053	20.000	ng/u1	0.00
20) Naphthalene-d8	10.439	136	458912	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.310	164	276511	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.062	188	528312	20.000	ng/u1	0.00
79) Chrysene-d12	21.262	240	359925	20.000	ng/u1	0.00
88) Perylene-d12	23.497	264	411236	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.234	96	15443	4.847	ng/uL	0.00
4) Pyridine-d5	3.628	84	257627	27.475	ng/u1	0.00
7) Phenol-d5	6.839	99	294742	26.532	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.016	67	199636	29.093	ng/u1	0.00
11) 2-Chlorophenol-d4	7.192	132	232759	28.511	ng/u1	0.00
15) 4-Methylphenol-d8	8.369	113	232331	26.630	ng/u1	0.00
21) Nitrobenzene-d5	8.822	128	119773	30.994	ng/u1	0.00
24) 2-Nitrophenol-d4	9.533	143	132477	31.801	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.063	165	212780	30.012	ng/u1	0.00
31) 4-Chloroaniline-d4	10.592	131	207448	17.617	ng/u1	0.00
46) Dimethylphthalate-d6	13.733	166	641657	29.582	ng/u1	0.00
49) Acenaphthylene-d8	13.998	160	745209	30.041	ng/u1	0.00
54) 4-Nitrophenol-d4	14.533	143	99684	22.161	ng/u1	0.00
60) Fluorene-d10	15.309	176	546882	30.003	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.445	200	92493	26.971	ng/u1	0.00
73) Anthracene-d10	17.162	188	834721	33.054	ng/u1	0.00
81) Pyrene-d10	19.468	212	842321	38.157	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.356	264	708960	32.799	ng/u1	0.00

Target Compounds Qvalue

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

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