

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM101424\
 Data File : BM048062.D
 Acq On : 15 Oct 2024 15:49
 Operator : RC/JU
 Sample : PB164081BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK081

Quant Time: Oct 15 23:30:30 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM092724.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Oct 10 05:10:14 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.781	152	298655	20.000	ng/u1	0.00
20) Naphthalene-d8	10.575	136	1216409	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.422	164	784684	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.157	188	1571801	20.000	ng/u1	0.00
79) Chrysene-d12	21.386	240	1431641	20.000	ng/u1	0.00
88) Perylene-d12	24.386	264	1584328	20.000	ng/u1	0.01

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.240	96	33429	3.617	ng/uL	0.00
4) Pyridine-d5	3.652	84	463233	17.119	ng/u1	0.00
7) Phenol-d5	6.957	99	541652	19.180	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.110	67	325562	16.246	ng/u1	0.00
11) 2-Chlorophenol-d4	7.316	132	470498	22.899	ng/u1	0.00
15) 4-Methylphenol-d8	8.492	113	425200	20.251	ng/u1	0.00
21) Nitrobenzene-d5	8.934	128	227396	23.479	ng/u1	0.00
24) 2-Nitrophenol-d4	9.657	143	253624	27.327	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.198	165	415900	22.432	ng/u1	0.00
31) 4-Chloroaniline-d4	10.710	131	614593	21.441	ng/u1	0.00
46) Dimethylphthalate-d6	13.827	166	1277632	22.458	ng/u1	0.00
49) Acenaphthylene-d8	14.110	160	1475368	21.873	ng/u1	0.00
54) 4-Nitrophenol-d4	14.633	143	202347	17.864	ng/u1	0.01
60) Fluorene-d10	15.410	176	1099411	22.311	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.533	200	185814	20.202	ng/u1	0.00
73) Anthracene-d10	17.257	188	1681491	21.756	ng/u1	0.00
81) Pyrene-d10	19.551	212	1925998	23.704	ng/u1	0.00
92) Benzo(a)pyrene-d12	24.180	264	1921455	23.375	ng/u1	0.01

Target Compounds	Qvalue

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

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