

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM092824\
 Data File : BM047688.D
 Acq On : 28 Sep 2024 12:44
 Operator : RC/JU
 Sample : PB163765BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK765

Quant Time: Sep 29 00:48:40 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM092724.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Sat Sep 28 02:30:36 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.816	152	193863	20.000	ng/u1	0.00
20) Naphthalene-d8	10.610	136	760972	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.445	164	443105	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.186	188	858061	20.000	ng/u1	0.00
79) Chrysene-d12	21.409	240	756789	20.000	ng/u1	0.00
88) Perylene-d12	24.415	264	881547	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.257	96	34453	5.743	ng/uL	0.00
4) Pyridine-d5	3.669	84	494296	28.141	ng/u1	0.00
7) Phenol-d5	6.975	99	511221	27.887	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.145	67	377596	29.028	ng/u1	0.00
11) 2-Chlorophenol-d4	7.345	132	394758	29.598	ng/u1	0.00
15) 4-Methylphenol-d8	8.516	113	379487	27.843	ng/u1	0.00
21) Nitrobenzene-d5	8.969	128	187219	30.900	ng/u1	0.00
24) 2-Nitrophenol-d4	9.686	143	181315	31.228	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.228	165	336297	28.994	ng/u1	0.00
31) 4-Chloroaniline-d4	10.739	131	528390	29.467	ng/u1	0.00
46) Dimethylphthalate-d6	13.857	166	1033403	32.168	ng/u1	0.00
49) Acenaphthylene-d8	14.139	160	1173184	30.801	ng/u1	0.00
54) 4-Nitrophenol-d4	14.633	143	160483	25.091	ng/u1	0.00
60) Fluorene-d10	15.439	176	878995	31.589	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.551	200	119747	23.848	ng/u1	0.00
73) Anthracene-d10	17.280	188	1325533	31.417	ng/u1	0.00
81) Pyrene-d10	19.568	212	1524539	35.494	ng/u1	0.00
92) Benzo(a)pyrene-d12	24.209	264	1432772	31.325	ng/u1	0.00

Target Compounds Qvalue

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

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