

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM110824\
 Data File : BM048533.D
 Acq On : 08 Nov 2024 16:56
 Operator : RC/JU
 Sample : PB164803BL
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK803

Quant Time: Nov 08 22:32:17 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM110724.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Nov 07 22:50:20 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.663	152	137708	20.000	ng/u1	0.00
20) Naphthalene-d8	10.451	136	558360	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.316	164	337902	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.057	188	683345	20.000	ng/u1	0.00
79) Chrysene-d12	21.286	240	672371	20.000	ng/u1	0.00
88) Perylene-d12	24.197	264	794523	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.128	96	19400	5.616	ng/uL	0.00
4) Pyridine-d5	3.546	84	249569	26.216	ng/u1	0.00
7) Phenol-d5	6.845	99	297265	28.357	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	6.998	67	176954	28.104	ng/u1	0.00
11) 2-Chlorophenol-d4	7.198	132	256288	29.514	ng/u1	0.00
15) 4-Methylphenol-d8	8.381	113	240991	29.237	ng/u1	0.00
21) Nitrobenzene-d5	8.822	128	119130	29.437	ng/u1	0.00
24) 2-Nitrophenol-d4	9.539	143	128123	29.765	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.081	165	215034	26.891	ng/u1	0.00
31) 4-Chloroaniline-d4	10.598	131	322171	28.364	ng/u1	0.00
46) Dimethylphthalate-d6	13.727	166	647450	30.016	ng/u1	0.00
49) Acenaphthylene-d8	14.004	160	780734	30.700	ng/u1	0.00
54) 4-Nitrophenol-d4	14.545	143	86770	21.659	ng/u1	0.01
60) Fluorene-d10	15.310	176	561253	29.809	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.445	200	73005	21.237	ng/u1	0.00
73) Anthracene-d10	17.157	188	889350	30.755	ng/u1	0.00
81) Pyrene-d10	19.451	212	1073340	31.809	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.997	264	1177853	31.729	ng/u1	0.00

Target Compounds	Qvalue

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

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