

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG080724\
 Data File : BG062298.D
 Acq On : 7 Aug 2024 11:57
 Operator : MA/JU
 Sample : PB162309BL
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SBLK309

Quant Time: Aug 08 06:16:17 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG080524.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Aug 07 01:46:39 2024
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.963	152	53215	20.000	ng/u1	0.00
20) Naphthalene-d8	10.753	136	249411	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.578	164	214801	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.321	188	474963	20.000	ng/u1	0.00
79) Chrysene-d12	21.562	240	448455	20.000	ng/u1	0.00
88) Perylene-d12	24.658	264	539576	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.422	96	6558	5.314	ng/uL	0.00
4) Pyridine-d5	3.827	84	102547	26.412	ng/u1	0.00
7) Phenol-d5	7.111	99	142795	28.113	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.287	67	86279	27.595	ng/u1	0.00
11) 2-Chlorophenol-d4	7.493	132	100540	27.628	ng/u1	0.00
15) 4-Methylphenol-d8	8.650	113	123782	29.258	ng/u1	0.00
21) Nitrobenzene-d5	9.114	128	54089	35.170	ng/u1	0.00
24) 2-Nitrophenol-d4	9.831	143	57761	39.925	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.366	165	128688	30.626	ng/u1	0.00
31) 4-Chloroaniline-d4	10.883	131	188475	30.926	ng/u1	0.00
46) Dimethylphthalate-d6	13.990	166	449819	27.142	ng/u1	0.00
49) Acenaphthylene-d8	14.272	160	510629	27.175	ng/u1	0.00
54) 4-Nitrophenol-d4	14.754	143	62929	27.122	ng/u1	0.00
60) Fluorene-d10	15.570	176	418216	28.114	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.676	200	44788	31.841	ng/u1	0.00
73) Anthracene-d10	17.421	188	639517	27.916	ng/u1	0.00
81) Pyrene-d10	19.700	212	769632	28.874	ng/u1	0.00
92) Benzo(a)pyrene-d12	24.447	264	797936	27.663	ng/u1	0.00

Target Compounds Qvalue

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

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