

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG112321\
 Data File : BG051189.D
 Acq On : 23 Nov 2021 18:37
 Operator : CG/JU
 Sample : PB140910BL
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SBLK910

Quant Time: Nov 24 06:55:30 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG112321.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Nov 24 06:04:50 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.196	152	22236	20.000	ng/u1	0.00
20) Naphthalene-d8	11.028	136	99079	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.830	164	72153	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.580	188	173113	20.000	ng/u1	0.00
79) Chrysene-d12	21.880	240	159789	20.000	ng/u1	0.00
88) Perylene-d12	25.282	264	156501	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.537	96	4095	6.400	ng/uL	0.00
4) Pyridine-d5	3.972	84	53467	28.476	ng/u1	0.00
7) Phenol-d5	7.356	99	64909	29.535	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.515	67	43579	31.574	ng/u1	0.00
11) 2-Chlorophenol-d4	7.732	132	49469	31.259	ng/u1	0.00
15) 4-Methylphenol-d8	8.913	113	53229	30.014	ng/u1	0.00
21) Nitrobenzene-d5	9.377	128	26330	31.481	ng/u1	0.00
24) 2-Nitrophenol-d4	10.100	143	30561	32.392	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.652	165	46108	28.804	ng/u1	0.00
31) 4-Chloroaniline-d4	11.164	131	73556	31.404	ng/u1	0.00
46) Dimethylphthalate-d6	14.225	166	189914	34.208	ng/u1	0.00
49) Acenaphthylene-d8	14.530	160	221608	31.655	ng/u1	0.00
54) 4-Nitrophenol-d4	15.047	143	25591	28.477	ng/u1	0.00
60) Fluorene-d10	15.823	176	161094	32.223	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.952	200	27787	26.012	ng/u1	0.00
73) Anthracene-d10	17.679	188	272460	32.908	ng/u1	0.00
81) Pyrene-d10	19.959	212	316646	32.750	ng/u1	0.00
92) Benzo(a)pyrene-d12	25.053	264	279195	33.404	ng/u1	0.01

Target Compounds Qvalue

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

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