

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG110321\
 Data File : BG050864.D
 Acq On : 4 Nov 2021 2:28
 Operator : CG/JU
 Sample : PB140455BL
 Misc :
 ALS Vial : 51 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SBLK455

Quant Time: Nov 07 08:05:27 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG110321.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Nov 02 14:49:05 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.226	152	30845	20.000	ng/u1	0.00
20) Naphthalene-d8	11.053	136	140802	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.854	164	100623	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.598	188	235099	20.000	ng/u1	0.00
79) Chrysene-d12	21.893	240	207987	20.000	ng/u1	-0.01
88) Perylene-d12	25.289	264	203309	20.000	ng/u1	-0.02

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.585	96	4965	5.195	ng/uL	0.00
4) Pyridine-d5	4.008	84	75002	26.234	ng/u1	0.00
7) Phenol-d5	7.369	99	85619	26.019	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.545	67	58859	27.690	ng/u1	0.00
11) 2-Chlorophenol-d4	7.756	132	61550	26.990	ng/u1	0.00
15) 4-Methylphenol-d8	8.926	113	67912	26.215	ng/u1	0.00
21) Nitrobenzene-d5	9.402	128	33768	28.221	ng/u1	0.00
24) 2-Nitrophenol-d4	10.130	143	36439	27.388	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.665	165	59030	26.339	ng/u1	0.00
31) 4-Chloroaniline-d4	11.182	131	93692	27.606	ng/u1	0.00
46) Dimethylphthalate-d6	14.243	166	234781	30.498	ng/u1	0.00
49) Acenaphthylene-d8	14.548	160	277955	28.980	ng/u1	0.00
54) 4-Nitrophenol-d4	15.036	143	36303	26.008	ng/u1	0.00
60) Fluorene-d10	15.841	176	198000	29.034	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.953	200	32478	22.786	ng/u1	0.00
73) Anthracene-d10	17.698	188	333913	30.041	ng/u1	0.00
81) Pyrene-d10	19.971	212	383592	28.555	ng/u1	0.00
92) Benzo(a)pyrene-d12	25.060	264	319855	28.459	ng/u1	-0.01

Target Compounds	Qvalue

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

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