

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG123021\
 Data File : BG051853.D
 Acq On : 30 Dec 2021 18:51
 Operator : CG/JU
 Sample : PB141709BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SBLK709

Quant Time: Dec 31 00:19:24 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG121421.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Dec 31 00:13:36 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.272	152	22429	20.000	ng/ul	0.00
20) Naphthalene-d8	11.109	136	96405	20.000	ng/ul	# 0.00
38) Acenaphthene-d10	14.904	164	71428	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.647	188	190224	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.965	240	180066	20.000	ng/ul	# 0.00
88) Perylene-d12	25.454	264	162788	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.590	96	4489	8.312	ng/uL	0.00
4) Pyridine-d5	4.025	84	61932	37.954	ng/ul	0.00
7) Phenol-d5	7.408	99	77993	38.946	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.579	67	49759	41.796	ng/ul	0.00
11) 2-Chlorophenol-d4	7.796	132	52150	37.442	ng/ul	0.00
15) 4-Methylphenol-d8	8.971	113	57738	37.299	ng/ul	0.00
21) Nitrobenzene-d5	9.441	128	27439	38.821	ng/ul	0.00
24) 2-Nitrophenol-d4	10.169	143	30286	39.197	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.716	165	55584	33.174	ng/ul	0.00
31) 4-Chloroaniline-d4	11.233	131	77154	34.974	ng/ul	0.00
46) Dimethylphthalate-d6	14.293	166	212602	37.187	ng/ul	0.00
49) Acenaphthylene-d8	14.599	160	254592	35.856	ng/ul	0.00
54) 4-Nitrophenol-d4	15.080	143	34176	36.760	ng/ul	0.00
60) Fluorene-d10	15.891	176	184679	36.055	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.997	200	36270	36.596	ng/ul	0.00
73) Anthracene-d10	17.747	188	330465	36.418	ng/ul	0.00
81) Pyrene-d10	20.027	212	414655	38.262	ng/ul	0.00
92) Benzo(a)pyrene-d12	25.220	264	324844	37.926	ng/ul	0.00

Target Compounds Qvalue

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

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