

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG122721\
 Data File : BG051796.D
 Acq On : 27 Dec 2021 10:29
 Operator : CG/JU
 Sample : PB141658BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SBLK658

Quant Time: Dec 27 11:05:16 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG121421.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Dec 23 13:37:07 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.276	152	30564	20.000	ng/u1	-0.01
20) Naphthalene-d8	11.113	136	127956	20.000	ng/u1	-0.01
38) Acenaphthene-d10	14.908	164	90099	20.000	ng/u1	-0.01
64) Phenanthrene-d10	17.657	188	199036	20.000	ng/u1	# 0.00
79) Chrysene-d12	21.975	240	180622	20.000	ng/u1	#-0.01
88) Perylene-d12	25.476	264	167446	20.000	ng/u1	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.600	96	5135	6.977	ng/uL	0.00
4) Pyridine-d5	4.023	84	79940	35.951	ng/u1	-0.02
7) Phenol-d5	7.407	99	95577	35.023	ng/u1	-0.01
9) Bis-(2-Chloroethyl)eth...	7.577	67	57471	35.426	ng/u1	-0.02
11) 2-Chlorophenol-d4	7.800	132	67365	35.493	ng/u1	-0.01
15) 4-Methylphenol-d8	8.969	113	73327	34.761	ng/u1	-0.01
21) Nitrobenzene-d5	9.445	128	34649	36.934	ng/u1	-0.01
24) 2-Nitrophenol-d4	10.179	143	37718	36.779	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.720	165	74585	33.538	ng/u1	-0.01
31) 4-Chloroaniline-d4	11.231	131	99972	34.143	ng/u1	-0.01
46) Dimethylphthalate-d6	14.297	166	237175	32.889	ng/u1	-0.02
49) Acenaphthylene-d8	14.603	160	305672	34.129	ng/u1	-0.01
54) 4-Nitrophenol-d4	15.073	143	28606	24.393	ng/u1	0.00
60) Fluorene-d10	15.895	176	209857	32.481	ng/u1	-0.01
65) 4,6-Dinitro-2-methylph...	16.001	200	20953	20.205	ng/u1	-0.01
73) Anthracene-d10	17.757	188	340456	35.858	ng/u1	0.00
81) Pyrene-d10	20.031	212	391720	36.034	ng/u1	-0.01
92) Benzo(a)pyrene-d12	25.230	264	329570	37.407	ng/u1	-0.02

Target Compounds	Qvalue

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

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