

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG011222\
 Data File : BG051963.D
 Acq On : 13 Jan 2022 9:24
 Operator : CG/JU
 Sample : PB141967BL
 Misc :
 ALS Vial : 28 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SBLK967

Quant Time: Jan 13 10:00:07 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG010622.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Jan 06 14:43:02 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.265	152	22288	20.000	ng/u1	0.01
20) Naphthalene-d8	11.102	136	92122	20.000	ng/u1	0.01
38) Acenaphthene-d10	14.897	164	69150	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.646	188	197660	20.000	ng/u1	# 0.00
79) Chrysene-d12	21.964	240	213395	20.000	ng/u1	# 0.00
88) Perylene-d12	25.459	264	200403	20.000	ng/u1	0.02

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.577	96	3756	6.022	ng/uL	0.00
4) Pyridine-d5	4.006	84	54232	27.970	ng/u1	0.00
7) Phenol-d5	7.401	99	63445	29.085	ng/u1	0.01
9) Bis-(2-Chloroethyl)eth...	7.566	67	38551	27.839	ng/u1	0.00
11) 2-Chlorophenol-d4	7.789	132	43311	29.988	ng/u1	0.01
15) 4-Methylphenol-d8	8.964	113	49186	29.006	ng/u1	0.01
21) Nitrobenzene-d5	9.434	128	22759	31.266	ng/u1	0.01
24) 2-Nitrophenol-d4	10.168	143	24513	30.448	ng/u1	0.02
28) 2,4-Dichlorophenol-d3	10.714	165	46091	29.290	ng/u1	0.01
31) 4-Chloroaniline-d4	11.225	131	63199	28.839	ng/u1	0.01
46) Dimethylphthalate-d6	14.292	166	180721	31.618	ng/u1	0.01
49) Acenaphthylene-d8	14.597	160	202099	29.271	ng/u1	0.01
54) 4-Nitrophenol-d4	15.073	143	31913	33.395	ng/u1	0.00
60) Fluorene-d10	15.890	176	151989	30.499	ng/u1	0.01
65) 4,6-Dinitro-2-methylph...	15.995	200	31951	25.624	ng/u1	0.01
73) Anthracene-d10	17.746	188	297677	31.747	ng/u1	0.00
81) Pyrene-d10	20.025	212	406105	33.067	ng/u1	0.01
92) Benzo(a)pyrene-d12	25.218	264	348737	33.055	ng/u1	0.02

Target Compounds Qvalue

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

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