

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG122721\
 Data File : BG051813.D
 Acq On : 27 Dec 2021 22:45
 Operator : CG/JU
 Sample : PB141643BL
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SBLK643

Manual Integrations
 APPROVED

Reviewed By :Jagrut Upadhyay 12/28/2021
 Supervised By :mohammad ahmed 12/31/2021

Quant Time: Dec 28 00:08:35 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	8.275	152	35878	20.000	ng/u1	-0.01
20) Naphthalene-d8	11.107	136	152850	20.000	ng/u1	-0.02
38) Acenaphthene-d10	14.908	164	104613	20.000	ng/u1	-0.01
64) Phenanthrene-d10	17.657	188	231849m	20.000	ng/u1	0.00
79) Chrysene-d12	21.974	240	228103m	20.000	ng/u1	-0.01
88) Perylene-d12	25.470	264	198365m	20.000	ng/u1	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.599	96	5063	5.861	ng/uL	0.00
4) Pyridine-d5	4.028	84	90760	34.771	ng/u1	-0.02
7) Phenol-d5	7.406	99	109698	34.244	ng/u1	-0.01
9) Bis-(2-Chloroethyl)eth...	7.576	67	65128	34.199	ng/u1	-0.02
11) 2-Chlorophenol-d4	7.800	132	76724	34.437	ng/u1	-0.01
15) 4-Methylphenol-d8	8.969	113	82373	33.266	ng/u1	-0.01
21) Nitrobenzene-d5	9.438	128	40476	36.119	ng/u1	-0.02
24) 2-Nitrophenol-d4	10.173	143	42630	34.799	ng/u1	-0.01
28) 2,4-Dichlorophenol-d3	10.719	165	80850	30.434	ng/u1	-0.01
31) 4-Chloroaniline-d4	11.230	131	105484	30.158	ng/u1	-0.01
46) Dimethylphthalate-d6	14.297	166	260189	31.074	ng/u1	-0.02
49) Acenaphthylene-d8	14.602	160	332319	31.957	ng/u1	-0.01
54) 4-Nitrophenol-d4	15.078	143	29156	21.412	ng/u1	0.00
60) Fluorene-d10	15.894	176	230188	30.684	ng/u1	-0.01
65) 4,6-Dinitro-2-methylph...	16.000	200	27492m	22.759	ng/u1	-0.01
73) Anthracene-d10	17.751	188	373148m	33.739	ng/u1	-0.01
81) Pyrene-d10	20.030	212	442225	32.212	ng/u1	-0.01
92) Benzo(a)pyrene-d12	25.229	264	359946m	34.487	ng/u1	-0.02

Target Compounds Qvalue

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

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