

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG083021\
 Data File : BG050090.D
 Acq On : 1 Sep 2021 9:35
 Operator : CG/JU
 Sample : PB138604BL
 Misc :
 ALS Vial : 64 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SBLK604

Quant Time: Sep 01 10:43:31 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG081921.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Aug 30 14:59:55 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.957	152	25766	20.000	ng/ul	0.00
20) Naphthalene-d8	10.776	136	105954	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.624	164	72669	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.373	188	154640	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.650	240	146656	20.000	ng/ul	0.00
88) Perylene-d12	24.875	264	146492	20.000	ng/ul	0.00

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.316	96	3354	6.712	ng/uL	0.00
4) Pyridine-d5	3.745	84	43574	28.952	ng/ul	0.00
7) Phenol-d5	7.128	99	66141	30.230	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.281	67	35760	29.419	ng/ul	0.00
11) 2-Chlorophenol-d4	7.493	132	51048	31.182	ng/ul	0.00
15) 4-Methylphenol-d8	8.685	113	52539	30.491	ng/ul	0.00
21) Nitrobenzene-d5	9.132	128	27348	33.293	ng/ul	0.00
24) 2-Nitrophenol-d4	9.860	143	31017	31.711	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.406	165	54976	31.512	ng/ul	0.00
31) 4-Chloroaniline-d4	10.917	131	71843	30.255	ng/ul	0.00
46) Dimethylphthalate-d6	14.019	166	173249	31.152	ng/ul	0.00
49) Acenaphthylene-d8	14.313	160	216580	33.214	ng/ul	0.00
54) 4-Nitrophenol-d4	14.853	143	20627	25.635	ng/ul	0.00
60) Fluorene-d10	15.617	176	149939	31.798	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.758	200	20139	20.243	ng/ul	0.00
73) Anthracene-d10	17.473	188	239522	34.493	ng/ul	0.00
81) Pyrene-d10	19.764	212	268277	33.851	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.652	264	248659	31.995	ng/ul	0.00

Target Compounds	Qvalue

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

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