

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG083021\
 Data File : BG050043.D
 Acq On : 30 Aug 2021 23:58
 Operator : CG/JU
 Sample : PB138632BL
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SBLK632

Quant Time: Aug 31 02:10:53 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	7.963	152	28044	20.000	ng/ul	0.00
20) Naphthalene-d8	10.783	136	118322	20.000	ng/ul	# 0.00
38) Acenaphthene-d10	14.625	164	84455	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.380	188	183010	20.000	ng/ul	0.00
79) Chrysene-d12	21.651	240	159875	20.000	ng/ul	0.00
88) Perylene-d12	24.876	264	163448	20.000	ng/ul	0.00

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.311	96	3512	6.458	ng/uL	0.00
4) Pyridine-d5	3.746	84	47103	28.754	ng/ul	0.00
7) Phenol-d5	7.135	99	77307	32.463	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.282	67	40998	30.988	ng/ul	0.00
11) 2-Chlorophenol-d4	7.493	132	59528	33.408	ng/ul	0.00
15) 4-Methylphenol-d8	8.686	113	59739	31.853	ng/ul	0.00
21) Nitrobenzene-d5	9.132	128	29613	32.282	ng/ul	0.00
24) 2-Nitrophenol-d4	9.861	143	35106	32.140	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.413	165	64237	32.972	ng/ul	0.00
31) 4-Chloroaniline-d4	10.918	131	83726	31.574	ng/ul	0.00
46) Dimethylphthalate-d6	14.020	166	201003	31.098	ng/ul	0.00
49) Acenaphthylene-d8	14.319	160	250644	33.073	ng/ul	0.00
54) 4-Nitrophenol-d4	14.854	143	26442	28.276	ng/ul	0.00
60) Fluorene-d10	15.618	176	176187	32.150	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.759	200	23182	19.689	ng/ul	0.00
73) Anthracene-d10	17.474	188	274807	33.439	ng/ul	0.00
81) Pyrene-d10	19.771	212	298706	34.574	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.658	264	278349	32.100	ng/ul	0.00

Target Compounds	Qvalue

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

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