

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG082621\
 Data File : BG049977.D
 Acq On : 26 Aug 2021 12:07
 Operator : CG/JU
 Sample : PB138646BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SBLK646

Quant Time: Aug 26 12:58:50 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG081921.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Aug 20 10:07:50 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.972	152	25912	20.000	ng/ul	-0.01
20) Naphthalene-d8	10.792	136	114971	20.000	ng/ul	-0.01
38) Acenaphthene-d10	14.633	164	84638	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.383	188	194990	20.000	ng/ul	-0.01
79) Chrysene-d12	21.659	240	145134	20.000	ng/ul	0.00
88) Perylene-d12	24.884	264	132354	20.000	ng/ul	#-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.331	96	3424	6.814	ng/uL	-0.01
4) Pyridine-d5	3.760	84	47977	31.698	ng/ul	-0.01
7) Phenol-d5	7.144	99	75439	34.285	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.290	67	43908	35.918	ng/ul	-0.02
11) 2-Chlorophenol-d4	7.502	132	59193	35.953	ng/ul	-0.01
15) 4-Methylphenol-d8	8.694	113	59207	34.167	ng/ul	0.00
21) Nitrobenzene-d5	9.147	128	31647	35.505	ng/ul	0.00
24) 2-Nitrophenol-d4	9.869	143	37249	35.096	ng/ul	-0.01
28) 2,4-Dichlorophenol-d3	10.421	165	61199	32.328	ng/ul	0.00
31) 4-Chloroaniline-d4	10.932	131	90987	35.312	ng/ul	0.00
46) Dimethylphthalate-d6	14.028	166	245604	37.917	ng/ul	-0.01
49) Acenaphthylene-d8	14.328	160	271704	35.775	ng/ul	0.00
54) 4-Nitrophenol-d4	14.862	143	30217	32.243	ng/ul	0.00
60) Fluorene-d10	15.626	176	198566	36.156	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.761	200	36577	29.157	ng/ul	0.00
73) Anthracene-d10	17.482	188	330468	37.742	ng/ul	-0.01
81) Pyrene-d10	19.773	212	369644	47.131	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.673	264	255310	36.360	ng/ul	0.00

Target Compounds	Qvalue

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

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