

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG082621\
 Data File : BG049983.D
 Acq On : 26 Aug 2021 20:14
 Operator : CG/JU
 Sample : PB138723BL
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SBLK723

Quant Time: Aug 27 02:01:07 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG081921.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Aug 26 15:41:36 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.968	152	26634	20.000	ng/u1	-0.01
20) Naphthalene-d8	10.788	136	110988	20.000	ng/u1	#-0.02
38) Acenaphthene-d10	14.630	164	81677	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.385	188	189151	20.000	ng/u1	0.00
79) Chrysene-d12	21.655	240	156675	20.000	ng/u1	0.00
88) Perylene-d12	24.875	264	138738	20.000	ng/u1	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.327	96	3373	6.531	ng/uL	-0.02
4) Pyridine-d5	3.756	84	48615	31.249	ng/u1	-0.02
7) Phenol-d5	7.140	99	76474	33.813	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.292	67	43996	35.015	ng/u1	-0.02
11) 2-Chlorophenol-d4	7.504	132	58023	34.287	ng/u1	0.00
15) 4-Methylphenol-d8	8.691	113	61803	34.698	ng/u1	0.00
21) Nitrobenzene-d5	9.143	128	32070	37.270	ng/u1	0.00
24) 2-Nitrophenol-d4	9.865	143	37560	36.659	ng/u1	-0.02
28) 2,4-Dichlorophenol-d3	10.424	165	60426	33.065	ng/u1	0.00
31) 4-Chloroaniline-d4	10.929	131	85312	34.298	ng/u1	0.00
46) Dimethylphthalate-d6	14.030	166	227196	36.346	ng/u1	0.00
49) Acenaphthylene-d8	14.324	160	261826	35.724	ng/u1	0.00
54) 4-Nitrophenol-d4	14.859	143	28693	31.727	ng/u1	0.00
60) Fluorene-d10	15.622	176	185892	35.075	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.757	200	30271	24.875	ng/u1	0.00
73) Anthracene-d10	17.479	188	312634	36.807	ng/u1	-0.02
81) Pyrene-d10	19.770	212	371965	43.933	ng/u1	0.00
92) Benzo(a)pyrene-d12	24.657	264	249166	33.852	ng/u1	-0.02

Target Compounds Qvalue

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

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