

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG093021\
 Data File : BG050350.D
 Acq On : 1 Oct 2021 9:33
 Operator : CG/JU
 Sample : PB139436BL
 Misc :
 ALS Vial : 32 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SBLK436

Quant Time: Oct 01 10:09:32 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG093021.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Sep 30 16:27:24 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.272	152	36622	20.00	ng/ul	0.00
20) Naphthalene-d8	11.098	136	157572	20.00	ng/ul	0.00
38) Acenaphthene-d10	14.887	164	113663	20.00	ng/ul	0.00
64) Phenanthrene-d10	17.631	188	269388	20.00	ng/ul	0.00
79) Chrysene-d12	21.932	240	248813	20.00	ng/ul	-0.02
88) Perylene-d12	25.358	264	241307	20.00	ng/ul	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.636	96	6216	6.24	ng/uL	0.00
4) Pyridine-d5	4.059	84	91280	30.73	ng/ul	-0.01
7) Phenol-d5	7.402	99	101251	31.64	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.584	67	67295	33.42	ng/ul	0.00
11) 2-Chlorophenol-d4	7.796	132	72065	32.14	ng/ul	0.00
15) 4-Methylphenol-d8	8.959	113	79485	32.82	ng/ul	0.00
21) Nitrobenzene-d5	9.441	128	37385	33.91	ng/ul	0.00
24) 2-Nitrophenol-d4	10.169	143	38591	32.62	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.698	165	70390	30.13	ng/ul	-0.01
31) 4-Chloroaniline-d4	11.221	131	105553	31.63	ng/ul	0.00
46) Dimethylphthalate-d6	14.282	166	254409	32.71	ng/ul	-0.01
49) Acenaphthylene-d8	14.582	160	307819	31.51	ng/ul	-0.01
54) 4-Nitrophenol-d4	15.046	143	48246	32.20	ng/ul	-0.01
60) Fluorene-d10	15.875	176	227810	32.51	ng/ul	-0.01
65) 4,6-Dinitro-2-methylph...	15.974	200	39171	29.15	ng/ul	-0.01
73) Anthracene-d10	17.731	188	389153	34.19	ng/ul	0.00
81) Pyrene-d10	20.005	212	457731	34.60	ng/ul	0.00
92) Benzo(a)pyrene-d12	25.122	264	380897	31.39	ng/ul	-0.02

Target Compounds Qvalue

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

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