

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG093021\
 Data File : BG050380.D
 Acq On : 2 Oct 2021 8:03
 Operator : CG/JU
 Sample : PB139462BL
 Misc :
 ALS Vial : 65 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SBLK462

Quant Time: Oct 04 01:04:54 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	8.267	152	48973	20.00	ng/ul	0.00
20) Naphthalene-d8	11.093	136	197711	20.00	ng/ul	0.00
38) Acenaphthene-d10	14.882	164	130233	20.00	ng/ul	0.00
64) Phenanthrene-d10	17.626	188	292330	20.00	ng/ul	0.00
79) Chrysene-d12	21.927	240	280420	20.00	ng/ul	-0.02
88) Perylene-d12	25.352	264	275791	20.00	ng/ul	-0.03
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.637	96	7338	5.50	ng/uL	0.00
4) Pyridine-d5	4.060	84	110623	27.85	ng/ul	0.00
7) Phenol-d5	7.397	99	123417	28.84	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.585	67	81825	30.39	ng/ul	0.00
11) 2-Chlorophenol-d4	7.791	132	89167	29.74	ng/ul	0.00
15) 4-Methylphenol-d8	8.954	113	92382	28.53	ng/ul	0.00
21) Nitrobenzene-d5	9.436	128	45615	32.98	ng/ul	0.00
24) 2-Nitrophenol-d4	10.164	143	48182	32.46	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.699	165	83585	28.52	ng/ul	0.00
31) 4-Chloroaniline-d4	11.222	131	121943	29.12	ng/ul	0.00
46) Dimethylphthalate-d6	14.277	166	278382	31.23	ng/ul	-0.02
49) Acenaphthylene-d8	14.583	160	333471	29.79	ng/ul	0.00
54) 4-Nitrophenol-d4	15.047	143	50117	29.19	ng/ul	0.00
60) Fluorene-d10	15.869	176	239581	29.84	ng/ul	-0.02
65) 4,6-Dinitro-2-methylph...	15.975	200	37092	25.44	ng/ul	0.00
73) Anthracene-d10	17.726	188	390424	31.61	ng/ul	0.00
81) Pyrene-d10	20.000	212	461215	30.93	ng/ul	0.00
92) Benzo(a)pyrene-d12	25.111	264	405824	29.26	ng/ul	-0.03

Target Compounds Qvalue

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

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