

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG092523\
 Data File : BG059177.D
 Acq On : 25 Sep 2023 10:21
 Operator : MA/JU
 Sample : PB155606BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SBLK606

Quant Time: Sep 25 22:36:34 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG091923.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Sep 25 22:33:28 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.311	152	52211	20.000	ng/ul	0.00
20) Naphthalene-d8	11.154	136	239478	20.000	ng/ul	# 0.00
38) Acenaphthene-d10	14.938	164	157268	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.688	188	385589	20.000	ng/ul	0.00
79) Chrysene-d12	22.012	240	305680	20.000	ng/ul	0.00
88) Perylene-d12	25.573	264	348248	20.000	ng/ul	0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.622	96	10042	7.528	ng/uL	0.00
4) Pyridine-d5	4.057	84	136240	34.815	ng/ul	0.00
7) Phenol-d5	7.435	99	187754	36.572	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.629	67	114841	36.452	ng/ul	0.00
11) 2-Chlorophenol-d4	7.829	132	137520	36.299	ng/ul	0.00
15) 4-Methylphenol-d8	8.998	113	137457	33.619	ng/ul	0.00
21) Nitrobenzene-d5	9.509	128	69208	37.452	ng/ul	0.00
24) 2-Nitrophenol-d4	10.226	143	74559	39.940	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.755	165	131247	37.033	ng/ul	0.00
31) 4-Chloroaniline-d4	11.301	131	191920	34.696	ng/ul	0.00
46) Dimethylphthalate-d6	14.333	166	469148	38.788	ng/ul	0.00
49) Acenaphthylene-d8	14.645	160	521425	35.406	ng/ul	0.01
54) 4-Nitrophenol-d4	15.120	143	46965	21.283	ng/ul	0.00
60) Fluorene-d10	15.925	176	407163	36.776	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	16.031	200	65701	29.971	ng/ul	0.00
73) Anthracene-d10	17.788	188	641398	36.558	ng/ul	0.00
81) Pyrene-d10	20.062	212	767080	42.933	ng/ul	0.00
92) Benzo(a)pyrene-d12	25.326	264	622254	36.593	ng/ul	0.01

Target Compounds Qvalue

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

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