

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG092823\
 Data File : BG059240.D
 Acq On : 29 Sep 2023 5:11
 Operator : MA/JU
 Sample : PB155823BL
 Misc :
 ALS Vial : 31 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SBLK823

Quant Time: Sep 29 06:16:46 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG092723.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Sep 28 04:33:25 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.230	152	36364	20.000	ng/u1	0.00
20) Naphthalene-d8	11.074	136	168495	20.000	ng/u1	# 0.00
38) Acenaphthene-d10	14.864	164	107122	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.613	188	245943	20.000	ng/u1	0.00
79) Chrysene-d12	21.914	240	231873	20.000	ng/u1	0.00
88) Perylene-d12	25.392	264	281529	20.000	ng/u1	-0.02

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.524	96	6799	7.057	ng/uL	0.00
4) Pyridine-d5	3.965	84	106127	37.751	ng/u1	0.00
7) Phenol-d5	7.355	99	142525	39.156	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.549	67	87561	37.656	ng/u1	0.00
11) 2-Chlorophenol-d4	7.748	132	100996	41.276	ng/u1	0.00
15) 4-Methylphenol-d8	8.923	113	104731	37.222	ng/u1	0.00
21) Nitrobenzene-d5	9.429	128	51285	43.133	ng/u1	0.00
24) 2-Nitrophenol-d4	10.146	143	56482	48.390	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.674	165	101778	42.446	ng/u1	0.00
31) 4-Chloroaniline-d4	11.215	131	153585	39.589	ng/u1	0.00
46) Dimethylphthalate-d6	14.258	166	345717	44.638	ng/u1	0.00
49) Acenaphthylene-d8	14.570	160	390096	40.128	ng/u1	0.00
54) 4-Nitrophenol-d4	15.052	143	58290	43.099	ng/u1	0.00
60) Fluorene-d10	15.851	176	308151	41.804	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.962	200	43006	34.018	ng/u1	0.00
73) Anthracene-d10	17.713	188	496337	44.290	ng/u1	0.00
81) Pyrene-d10	19.987	212	583835	44.962	ng/u1	0.00
92) Benzo(a)pyrene-d12	25.146	264	606816	44.854	ng/u1	-0.02

Target Compounds	Qvalue

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

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