

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG100723\
 Data File : BG059377.D
 Acq On : 8 Oct 2023 15:53
 Operator : MA/JU
 Sample : PB156054BL
 Misc :
 ALS Vial : 35 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SBLK054

Quant Time: Oct 09 01:42:04 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG100723.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Sat Oct 07 22:45:22 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.230	152	45894	20.000	ng/u1	0.00
20) Naphthalene-d8	11.074	136	223816	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.864	164	154809	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.608	188	369622	20.000	ng/u1	# 0.00
79) Chrysene-d12	21.914	240	341374	20.000	ng/u1	0.00
88) Perylene-d12	25.399	264	400379	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.524	96	9263	7.196	ng/uL	0.00
4) Pyridine-d5	3.959	84	147858	39.908	ng/u1	0.00
7) Phenol-d5	7.355	99	185460	36.598	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.549	67	109062	38.077	ng/u1	0.00
11) 2-Chlorophenol-d4	7.743	132	140117	39.535	ng/u1	0.00
15) 4-Methylphenol-d8	8.918	113	145617	36.881	ng/u1	0.00
21) Nitrobenzene-d5	9.429	128	72921	38.273	ng/u1	0.00
24) 2-Nitrophenol-d4	10.146	143	83959	40.714	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.675	165	146130	39.260	ng/u1	0.00
31) 4-Chloroaniline-d4	11.215	131	226288	38.274	ng/u1	0.00
46) Dimethylphthalate-d6	14.259	166	509336	40.159	ng/u1	0.00
49) Acenaphthylene-d8	14.564	160	607987	40.282	ng/u1	0.00
54) 4-Nitrophenol-d4	15.052	143	85500	35.974	ng/u1	0.00
60) Fluorene-d10	15.851	176	477008	39.402	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.963	200	88557	36.288	ng/u1	0.00
73) Anthracene-d10	17.713	188	751107	41.425	ng/u1	0.00
81) Pyrene-d10	19.987	212	849938	41.099	ng/u1	0.00
92) Benzo(a)pyrene-d12	25.158	264	872600	41.417	ng/u1	0.00

Target Compounds Qvalue

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

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