

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG100523\
 Data File : BG059289.D
 Acq On : 5 Oct 2023 9:54
 Operator : MA/JU
 Sample : PB155951BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SBLK951

Quant Time: Oct 05 22:47:04 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	8.224	152	35803	20.000	ng/u1	0.00
20) Naphthalene-d8	11.068	136	172245	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.864	164	111300	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.613	188	247474	20.000	ng/u1	0.00
79) Chrysene-d12	21.908	240	223971	20.000	ng/u1	-0.01
88) Perylene-d12	25.392	264	266049	20.000	ng/u1	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.524	96	6209	6.545	ng/uL	0.00
4) Pyridine-d5	3.959	84	85920	31.042	ng/u1	0.00
7) Phenol-d5	7.355	99	111022	30.979	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.549	67	69539	30.374	ng/u1	0.00
11) 2-Chlorophenol-d4	7.748	132	84566	35.103	ng/u1	0.00
15) 4-Methylphenol-d8	8.923	113	85855	30.991	ng/u1	0.00
21) Nitrobenzene-d5	9.429	128	41491	34.136	ng/u1	0.00
24) 2-Nitrophenol-d4	10.146	143	46220	38.736	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.668	165	84100	34.310	ng/u1	0.00
31) 4-Chloroaniline-d4	11.215	131	125207	31.571	ng/u1	0.00
46) Dimethylphthalate-d6	14.258	166	280765	34.891	ng/u1	0.00
49) Acenaphthylene-d8	14.564	160	335016	33.169	ng/u1	0.00
54) 4-Nitrophenol-d4	15.052	143	43054	30.639	ng/u1	0.00
60) Fluorene-d10	15.851	176	258165	33.708	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.962	200	31407	24.690	ng/u1	0.00
73) Anthracene-d10	17.713	188	391861	34.751	ng/u1	0.00
81) Pyrene-d10	19.987	212	443937	35.394	ng/u1	0.00
92) Benzo(a)pyrene-d12	25.146	264	443898	34.720	ng/u1	-0.02

Target Compounds Qvalue

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

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