

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091324\
 Data File : BG062873.D
 Acq On : 16 Sep 2024 9:51
 Operator : JU/RC
 Sample : PB163398BL
 Misc :
 ALS Vial : 64 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SBLK398

Quant Time: Sep 16 23:32:47 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG090924.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Sep 10 01:28:17 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.900	152	31471	20.000	ng/u1	0.00
20) Naphthalene-d8	10.691	136	125627	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.528	164	111042	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.271	188	305608	20.000	ng/u1	# 0.00
79) Chrysene-d12	21.519	240	349408	20.000	ng/u1	# 0.00
88) Perylene-d12	24.581	264	397713	20.000	ng/u1	-0.01

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.364	96	3971	5.704	ng/uL	0.00
4) Pyridine-d5	3.764	84	64546	29.126	ng/u1	0.00
7) Phenol-d5	7.072	99	78940	28.605	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.230	67	44234	26.735	ng/u1	0.00
11) 2-Chlorophenol-d4	7.436	132	60990	30.341	ng/u1	0.00
15) 4-Methylphenol-d8	8.605	113	70021	30.448	ng/u1	0.00
21) Nitrobenzene-d5	9.052	128	32914	36.658	ng/u1	0.00
24) 2-Nitrophenol-d4	9.774	143	38791	39.504	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.315	165	79036	36.959	ng/u1	0.00
31) 4-Chloroaniline-d4	10.826	131	93267	31.729	ng/u1	0.00
46) Dimethylphthalate-d6	13.934	166	279933	32.737	ng/u1	0.00
49) Acenaphthylene-d8	14.222	160	307469	32.376	ng/u1	0.00
54) 4-Nitrophenol-d4	14.727	143	41346	28.463	ng/u1	0.00
60) Fluorene-d10	15.521	176	264251	34.294	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.638	200	52514	29.873	ng/u1	0.00
73) Anthracene-d10	17.371	188	503152	34.886	ng/u1	0.00
81) Pyrene-d10	19.663	212	662049	33.671	ng/u1	0.00
92) Benzo(a)pyrene-d12	24.375	264	714936	34.044	ng/u1	0.00

Target Compounds	Qvalue

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

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