

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG111324\
 Data File : BG063371.D
 Acq On : 14 Nov 2024 1:39
 Operator : RC/JU
 Sample : P4802-02
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :

Quant Time: Nov 14 03:33:59 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG110624.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Nov 06 15:45:52 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.828	152	99947	20.000	ng/ul	-0.01
20) Naphthalene-d8	10.619	136	428175	20.000	ng/ul	# 0.00
38) Acenaphthene-d10	14.468	164	360873	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.217	188	940084	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.471	240	995807	20.000	ng/ul	# 0.00
88) Perylene-d12	24.503	264	1073392	20.000	ng/ul	-0.01

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.298	96	9311	4.212	ng/uL	0.00
4) Pyridine-d5	3.710	84	26212	3.588	ng/ul	0.00
7) Phenol-d5	7.006	99	96456	10.695	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.159	67	149257	28.215	ng/ul	-0.01
11) 2-Chlorophenol-d4	7.364	132	168092	28.715	ng/ul	-0.01
15) 4-Methylphenol-d8	8.539	113	175230	22.923	ng/ul	0.00
21) Nitrobenzene-d5	8.986	128	94080	31.254	ng/ul	0.00
24) 2-Nitrophenol-d4	9.703	143	111167	28.574	ng/ul	-0.01
28) 2,4-Dichlorophenol-d3	10.249	165	220359	28.862	ng/ul	0.00
31) 4-Chloroaniline-d4	10.754	131	170776	17.173	ng/ul	0.00
46) Dimethylphthalate-d6	13.874	166	845165	29.899	ng/ul	0.00
49) Acenaphthylene-d8	14.162	160	867451	31.353	ng/ul	0.00
54) 4-Nitrophenol-d4	14.691	143	32479	8.392	ng/ul	0.01
60) Fluorene-d10	15.466	176	777867	32.606	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.590	200	182770	30.935	ng/ul	0.00
73) Anthracene-d10	17.317	188	1505859	37.270	ng/ul	0.00
81) Pyrene-d10	19.620	212	1983194	39.722	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.297	264	1882011	36.315	ng/ul	0.00

Target Compounds Qvalue

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

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