

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG111324\
 Data File : BG063375.D
 Acq On : 14 Nov 2024 4:48
 Operator : RC/JU
 Sample : PB164924BL
 Misc :
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :

Quant Time: Nov 14 04:27:11 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	7.831	152	96321	20.000	ng/ul	0.00
20) Naphthalene-d8	10.622	136	417905	20.000	ng/ul	# 0.00
38) Acenaphthene-d10	14.471	164	361940	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.220	188	929108	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.474	240	979062	20.000	ng/ul	# 0.00
88) Perylene-d12	24.506	264	1044693	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.348	96	121	0.057	ng/uL	0.04
4) Pyridine-d5	3.701	84	207549	29.479	ng/ul	0.00
7) Phenol-d5	7.003	99	242911	27.948	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.162	67	159679	31.321	ng/ul	0.00
11) 2-Chlorophenol-d4	7.367	132	180743	32.039	ng/ul	0.00
15) 4-Methylphenol-d8	8.536	113	220148	29.884	ng/ul	0.00
21) Nitrobenzene-d5	8.983	128	95956	32.660	ng/ul	0.00
24) 2-Nitrophenol-d4	9.706	143	126295	33.260	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.246	165	223219	29.955	ng/ul	0.00
31) 4-Chloroaniline-d4	10.757	131	275742	28.409	ng/ul	0.00
46) Dimethylphthalate-d6	13.877	166	849115	29.950	ng/ul	0.00
49) Acenaphthylene-d8	14.159	160	922612	33.248	ng/ul	0.00
54) 4-Nitrophenol-d4	14.682	143	116428	29.994	ng/ul	0.00
60) Fluorene-d10	15.464	176	778610	32.541	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.593	200	171411	29.355	ng/ul	0.00
73) Anthracene-d10	17.320	188	1384436	34.670	ng/ul	0.00
81) Pyrene-d10	19.618	212	1826370	37.206	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.300	264	1750856	34.712	ng/ul	0.00

Target Compounds Qvalue

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

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