

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG122624\
 Data File : BG063892.D
 Acq On : 28 Dec 2024 7:12
 Operator : RC/JU
 Sample : PB165853BL
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SBLK853

Quant Time: Dec 29 23:21:23 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG121124.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Dec 24 00:44:39 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.795	152	116770	20.000	ng/ul	0.00
20) Naphthalene-d8	10.586	136	452520	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.440	164	336848	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.196	188	809986	20.000	ng/ul	0.00
79) Chrysene-d12	21.449	240	866114	20.000	ng/ul	0.01
88) Perylene-d12	24.469	264	886927	20.000	ng/ul	0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.259	96	12459	4.181	ng/uL	0.00
4) Pyridine-d5	3.664	84	170769	18.167	ng/ul	0.00
7) Phenol-d5	6.984	99	183146	17.030	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.131	67	119173	16.044	ng/ul	0.00
11) 2-Chlorophenol-d4	7.337	132	126169	18.121	ng/ul	0.00
15) 4-Methylphenol-d8	8.518	113	137935	16.520	ng/ul	0.00
21) Nitrobenzene-d5	8.952	128	61222	18.948	ng/ul	0.00
24) 2-Nitrophenol-d4	9.675	143	68044	18.787	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.221	165	128987	18.395	ng/ul	0.00
31) 4-Chloroaniline-d4	10.727	131	155486	17.463	ng/ul	0.00
46) Dimethylphthalate-d6	13.847	166	422539	18.435	ng/ul	0.00
49) Acenaphthylene-d8	14.135	160	492997	18.868	ng/ul	0.00
54) 4-Nitrophenol-d4	14.681	143	40295	12.408	ng/ul	0.01
60) Fluorene-d10	15.439	176	380625	18.782	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.580	200	29341	6.977	ng/ul	0.01
73) Anthracene-d10	17.296	188	688471	19.770	ng/ul	0.00
81) Pyrene-d10	19.593	212	857408	19.024	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.258	264	807782	19.394	ng/ul	0.01

Target Compounds	Qvalue

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

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