

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG111124\
 Data File : BG063338.D
 Acq On : 12 Nov 2024 23:44
 Operator : RC/JU
 Sample : PB164897BL
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SBLK897

Quant Time: Nov 12 23:24:13 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	7.829	152	99707	20.000	ng/u1	-0.01
20) Naphthalene-d8	10.620	136	472394	20.000	ng/u1	# 0.00
38) Acenaphthene-d10	14.469	164	430666	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.218	188	1022516	20.000	ng/u1	# 0.00
79) Chrysene-d12	21.472	240	1054179	20.000	ng/u1	# 0.00
88) Perylene-d12	24.510	264	1121110	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.294	96	12523	5.679	ng/uL	-0.01
4) Pyridine-d5	3.699	84	201471	27.644	ng/u1	-0.01
7) Phenol-d5	7.007	99	290655	32.306	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.165	67	174049	32.981	ng/u1	0.00
11) 2-Chlorophenol-d4	7.365	132	209548	35.883	ng/u1	-0.01
15) 4-Methylphenol-d8	8.540	113	259518	34.031	ng/u1	0.00
21) Nitrobenzene-d5	8.987	128	112078	33.748	ng/u1	0.00
24) 2-Nitrophenol-d4	9.710	143	142650	33.234	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.244	165	279513	33.183	ng/u1	0.00
31) 4-Chloroaniline-d4	10.755	131	336464	30.667	ng/u1	0.00
46) Dimethylphthalate-d6	13.875	166	992286	29.415	ng/u1	0.00
49) Acenaphthylene-d8	14.163	160	1099639	33.304	ng/u1	0.00
54) 4-Nitrophenol-d4	14.680	143	118902	25.743	ng/u1	0.00
60) Fluorene-d10	15.467	176	918864	32.274	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.591	200	170897	26.594	ng/u1	0.00
73) Anthracene-d10	17.318	188	1517487	34.530	ng/u1	0.00
81) Pyrene-d10	19.621	212	1886873	35.700	ng/u1	0.00
92) Benzo(a)pyrene-d12	24.304	264	1861847	34.397	ng/u1	0.00

Target Compounds Qvalue

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

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