

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG111124\
 Data File : BG063339.D
 Acq On : 12 Nov 2024 00:22
 Operator : RC/JU
 Sample : PB164899BL
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SBLK899

Manual Integrations
 APPROVED

Reviewed By :Jagrut Upadhyay 11/13/2024
 Supervised By :mohammad ahmed 11/14/2024

Quant Time: Nov 13 00:10:32 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	7.830	152	103752	20.000	ng/u1	-0.01
20) Naphthalene-d8	10.621	136	496265	20.000	ng/u1	# 0.00
38) Acenaphthene-d10	14.469	164	453467	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.225	188	1048190	20.000	ng/u1	0.00
79) Chrysene-d12	21.473	240	1096735m	20.000	ng/u1	0.00
88) Perylene-d12	24.510	264	1172257	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.300	96	13185	5.746	ng/uL	0.00
4) Pyridine-d5	3.699	84	200517	26.440	ng/u1	-0.01
7) Phenol-d5	7.007	99	271969	29.050	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.160	67	168273	30.643	ng/u1	-0.01
11) 2-Chlorophenol-d4	7.366	132	193017	31.764	ng/u1	-0.01
15) 4-Methylphenol-d8	8.541	113	248380	31.301	ng/u1	0.00
21) Nitrobenzene-d5	8.987	128	107485	30.808	ng/u1	0.00
24) 2-Nitrophenol-d4	9.704	143	143570	31.839	ng/u1	-0.01
28) 2,4-Dichlorophenol-d3	10.244	165	274280	30.995	ng/u1	0.00
31) 4-Chloroaniline-d4	10.756	131	330773	28.698	ng/u1	0.00
46) Dimethylphthalate-d6	13.876	166	961638	27.073	ng/u1	0.00
49) Acenaphthylene-d8	14.163	160	1068823	30.743	ng/u1	0.00
54) 4-Nitrophenol-d4	14.680	143	118356	24.337	ng/u1	0.00
60) Fluorene-d10	15.468	176	894446	29.837	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.591	200	165780	25.165	ng/u1	0.00
73) Anthracene-d10	17.319	188	1460744	32.425	ng/u1	0.00
81) Pyrene-d10	19.622	212	1815734	33.021	ng/u1	0.00
92) Benzo(a)pyrene-d12	24.299	264	1845241	32.603	ng/u1	0.00

Target Compounds Qvalue

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

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