

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG111524\
 Data File : BG063452.D
 Acq On : 16 Nov 2024 10:14
 Operator : RC/JU
 Sample : PB164999BL
 Misc :
 ALS Vial : 37 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SBLK999

Quant Time: Nov 18 01:36:12 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	7.826	152	82642	20.000	ng/u1	-0.02
20) Naphthalene-d8	10.617	136	360842	20.000	ng/u1	# 0.00
38) Acenaphthene-d10	14.471	164	292107	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.226	188	741076	20.000	ng/u1	# 0.00
79) Chrysene-d12	21.480	240	795808	20.000	ng/u1	# 0.00
88) Perylene-d12	24.518	264	825223	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.302	96	11640	6.369	ng/uL	0.00
4) Pyridine-d5	3.701	84	187936	31.111	ng/u1	0.00
7) Phenol-d5	7.003	99	255694	34.288	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.162	67	161319	36.881	ng/u1	0.00
11) 2-Chlorophenol-d4	7.362	132	184956	38.212	ng/u1	-0.02
15) 4-Methylphenol-d8	8.537	113	205987	32.589	ng/u1	0.00
21) Nitrobenzene-d5	8.983	128	91215	35.956	ng/u1	0.00
24) 2-Nitrophenol-d4	9.706	143	114785	35.009	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.246	165	219027	34.040	ng/u1	0.00
31) 4-Chloroaniline-d4	10.758	131	276294	32.968	ng/u1	0.00
46) Dimethylphthalate-d6	13.877	166	755294	33.010	ng/u1	0.00
49) Acenaphthylene-d8	14.165	160	825036	36.840	ng/u1	0.00
54) 4-Nitrophenol-d4	14.694	143	91996	29.366	ng/u1	0.01
60) Fluorene-d10	15.470	176	696930	36.090	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.593	200	131941	28.329	ng/u1	0.00
73) Anthracene-d10	17.326	188	1247829	39.178	ng/u1	0.00
81) Pyrene-d10	19.624	212	1575056	39.475	ng/u1	0.00
92) Benzo(a)pyrene-d12	24.312	264	1552025	38.954	ng/u1	0.00

Target Compounds Qvalue

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

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