

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG111124\
 Data File : BG063336.D
 Acq On : 12 Nov 2024 21:51
 Operator : RC/JU
 Sample : PB164901BL
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SBLK901

Quant Time: Nov 12 22:45:37 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	7.832	152	89292	20.000	ng/u1	0.00
20) Naphthalene-d8	10.623	136	431792	20.000	ng/u1	# 0.00
38) Acenaphthene-d10	14.471	164	394011	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.221	188	928212	20.000	ng/u1	0.00
79) Chrysene-d12	21.475	240	961212	20.000	ng/u1	# 0.00
88) Perylene-d12	24.506	264	1019887	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.296	96	12466	6.313	ng/uL	0.00
4) Pyridine-d5	3.701	84	184869	28.324	ng/u1	0.00
7) Phenol-d5	7.003	99	265939	33.006	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.162	67	155594	32.923	ng/u1	0.00
11) 2-Chlorophenol-d4	7.368	132	178400	34.113	ng/u1	0.00
15) 4-Methylphenol-d8	8.543	113	233299	34.162	ng/u1	0.00
21) Nitrobenzene-d5	8.989	128	102222	33.674	ng/u1	0.00
24) 2-Nitrophenol-d4	9.706	143	132814	33.852	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.247	165	257232	33.409	ng/u1	0.00
31) 4-Chloroaniline-d4	10.758	131	310136	30.925	ng/u1	0.00
46) Dimethylphthalate-d6	13.878	166	920602	29.829	ng/u1	0.00
49) Acenaphthylene-d8	14.160	160	1021068	33.801	ng/u1	0.00
54) 4-Nitrophenol-d4	14.683	143	111685	26.430	ng/u1	0.00
60) Fluorene-d10	15.464	176	832517	31.962	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.593	200	152270	26.102	ng/u1	0.00
73) Anthracene-d10	17.321	188	1409165	35.323	ng/u1	0.00
81) Pyrene-d10	19.618	212	1741679	36.140	ng/u1	0.00
92) Benzo(a)pyrene-d12	24.301	264	1709193	34.710	ng/u1	0.00

Target Compounds Qvalue

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

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