

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG110824\
 Data File : BG063233.D
 Acq On : 8 Nov 2024 17:59
 Operator : RC/JU
 Sample : PB164573BL
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SBLK573

Quant Time: Nov 08 17:39:14 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	7.838	152	96762	20.000	ng/u1	0.00
20) Naphthalene-d8	10.629	136	458550	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.472	164	407829	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.227	188	1005060	20.000	ng/u1	0.00
79) Chrysene-d12	21.481	240	1027104	20.000	ng/u1	0.00
88) Perylene-d12	24.519	264	1078069	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.296	96	13869	6.481	ng/uL	0.00
4) Pyridine-d5	3.708	84	216206	30.568	ng/u1	0.00
7) Phenol-d5	7.010	99	271329	31.076	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.168	67	167263	32.659	ng/u1	0.00
11) 2-Chlorophenol-d4	7.368	132	194752	34.364	ng/u1	0.00
15) 4-Methylphenol-d8	8.543	113	233145	31.504	ng/u1	0.00
21) Nitrobenzene-d5	8.990	128	104947	32.554	ng/u1	0.00
24) 2-Nitrophenol-d4	9.712	143	137531	33.009	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.253	165	265281	32.444	ng/u1	0.00
31) 4-Chloroaniline-d4	10.764	131	317072	29.772	ng/u1	0.00
46) Dimethylphthalate-d6	13.878	166	925012	28.956	ng/u1	0.00
49) Acenaphthylene-d8	14.166	160	1040796	33.287	ng/u1	0.00
54) 4-Nitrophenol-d4	14.683	143	121902	27.871	ng/u1	0.00
60) Fluorene-d10	15.470	176	849701	31.516	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.594	200	175762	27.826	ng/u1	0.00
73) Anthracene-d10	17.321	188	1469783	34.026	ng/u1	0.00
81) Pyrene-d10	19.624	212	1833152	35.598	ng/u1	0.00
92) Benzo(a)pyrene-d12	24.313	264	1786365	34.320	ng/u1	0.00

Target Compounds Qvalue

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

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