

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG110824\
 Data File : BG063264.D
 Acq On : 9 Nov 2024 14:52
 Operator : RC/JU
 Sample : PB164616BL
 Misc :
 ALS Vial : 42 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SBLK616

Quant Time: Nov 10 01:11:04 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	7.835	152	137601	20.000	ng/u1	0.00
20) Naphthalene-d8	10.625	136	619346	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.474	164	537839	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.224	188	1238558	20.000	ng/u1	0.00
79) Chrysene-d12	21.477	240	1263072	20.000	ng/u1	# 0.00
88) Perylene-d12	24.509	264	1369310	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.305	96	16445	5.404	ng/uL	0.00
4) Pyridine-d5	3.704	84	246676	24.525	ng/u1	0.00
7) Phenol-d5	7.006	99	327526	26.379	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.165	67	194085	26.649	ng/u1	0.00
11) 2-Chlorophenol-d4	7.370	132	226902	28.155	ng/u1	0.00
15) 4-Methylphenol-d8	8.540	113	282283	26.823	ng/u1	0.00
21) Nitrobenzene-d5	8.992	128	132056	30.328	ng/u1	0.00
24) 2-Nitrophenol-d4	9.709	143	156536	27.816	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.249	165	313615	28.397	ng/u1	0.00
31) 4-Chloroaniline-d4	10.760	131	357876	24.879	ng/u1	0.00
46) Dimethylphthalate-d6	13.880	166	1055671	25.058	ng/u1	0.00
49) Acenaphthylene-d8	14.162	160	1183147	28.693	ng/u1	0.00
54) 4-Nitrophenol-d4	14.679	143	133735	23.185	ng/u1	0.00
60) Fluorene-d10	15.467	176	964039	27.114	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.590	200	160432	20.610	ng/u1	0.00
73) Anthracene-d10	17.323	188	1573952	29.568	ng/u1	0.00
81) Pyrene-d10	19.621	212	1983011	31.314	ng/u1	0.00
92) Benzo(a)pyrene-d12	24.303	264	1991388	30.121	ng/u1	0.00

Target Compounds Qvalue

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

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