

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG102721\
 Data File : BG050747.D
 Acq On : 27 Oct 2021 17:46
 Operator : CG/JU
 Sample : PB140133BL
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SBLK133

Quant Time: Oct 27 18:22:12 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG102721.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Oct 26 18:31:25 2021
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.248	152	30351	20.000	ng/u1	0.00
20) Naphthalene-d8	11.074	136	140629	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.869	164	97967	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.619	188	218023	20.000	ng/u1	0.00
79) Chrysene-d12	21.926	240	189902	20.000	ng/u1	# 0.00
88) Perylene-d12	25.345	264	183248	20.000	ng/u1	-0.02

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.600	96	4707	6.111	ng/uL	0.00
4) Pyridine-d5	4.023	84	66132	27.110	ng/u1	0.00
7) Phenol-d5	7.384	99	82719	29.329	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.560	67	54745	30.169	ng/u1	0.00
11) 2-Chlorophenol-d4	7.772	132	57857	29.204	ng/u1	0.00
15) 4-Methylphenol-d8	8.941	113	66170	30.386	ng/u1	0.00
21) Nitrobenzene-d5	9.417	128	31834	29.657	ng/u1	0.00
24) 2-Nitrophenol-d4	10.146	143	35194	29.075	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.686	165	56065	26.374	ng/u1	0.00
31) 4-Chloroaniline-d4	11.203	131	87381	28.008	ng/u1	0.00
46) Dimethylphthalate-d6	14.264	166	223714	32.335	ng/u1	0.00
49) Acenaphthylene-d8	14.570	160	249993	29.095	ng/u1	0.00
54) 4-Nitrophenol-d4	15.052	143	35686	28.500	ng/u1	0.00
60) Fluorene-d10	15.862	176	183976	30.942	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.974	200	32378	25.378	ng/u1	0.00
73) Anthracene-d10	17.719	188	310040	33.167	ng/u1	0.00
81) Pyrene-d10	19.993	212	357239	33.010	ng/u1	0.00
92) Benzo(a)pyrene-d12	25.110	264	294684	31.255	ng/u1	-0.01

Target Compounds	Qvalue

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

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