

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG110321\
 Data File : BG050818.D
 Acq On : 2 Nov 2021 15:36
 Operator : CG/JU
 Sample : PB140311BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SBLK311

Quant Time: Nov 02 16:27:25 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG110321.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Nov 02 14:49:05 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.231	152	38305	20.000	ng/u1	0.00
20) Naphthalene-d8	11.057	136	186345	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.853	164	132177	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.603	188	315071	20.000	ng/u1	0.00
79) Chrysene-d12	21.898	240	278631	20.000	ng/u1	0.00
88) Perylene-d12	25.300	264	245955	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.590	96	4950	4.171	ng/uL	0.00
4) Pyridine-d5	4.013	84	70458	19.845	ng/u1	0.00
7) Phenol-d5	7.374	99	81823	20.023	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.544	67	57338	21.721	ng/u1	0.00
11) 2-Chlorophenol-d4	7.755	132	59158	20.889	ng/u1	0.00
15) 4-Methylphenol-d8	8.931	113	64302	19.988	ng/u1	0.00
21) Nitrobenzene-d5	9.406	128	33177	20.951	ng/u1	0.00
24) 2-Nitrophenol-d4	10.129	143	36198	20.558	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.670	165	53621	18.078	ng/u1	0.00
31) 4-Chloroaniline-d4	11.187	131	89870	20.008	ng/u1	0.00
46) Dimethylphthalate-d6	14.248	166	227772	22.524	ng/u1	0.00
49) Acenaphthylene-d8	14.553	160	261602	20.764	ng/u1	0.00
54) 4-Nitrophenol-d4	15.041	143	35573	19.401	ng/u1	0.00
60) Fluorene-d10	15.846	176	190440	21.259	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.958	200	32320	16.920	ng/u1	0.00
73) Anthracene-d10	17.697	188	323937	21.746	ng/u1	0.00
81) Pyrene-d10	19.976	212	380929	21.167	ng/u1	0.00
92) Benzo(a)pyrene-d12	25.065	264	306599	22.549	ng/u1	0.00

Target Compounds Qvalue

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

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