

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG110821\
 Data File : BG050877.D
 Acq On : 8 Nov 2021 12:49
 Operator : CG/JU
 Sample : PB140507BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SBLK507

Quant Time: Nov 08 12:35:11 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	8.232	152	35701	20.000	ng/u1	0.00
20) Naphthalene-d8	11.058	136	159352	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.854	164	106073	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.597	188	235497	20.000	ng/u1	0.00
79) Chrysene-d12	21.898	240	208402	20.000	ng/u1	0.00
88) Perylene-d12	25.294	264	202320	20.000	ng/u1	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.590	96	5414	4.895	ng/uL	0.00
4) Pyridine-d5	4.013	84	79822	24.122	ng/u1	0.00
7) Phenol-d5	7.374	99	88688	23.286	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.544	67	61461	24.981	ng/u1	0.00
11) 2-Chlorophenol-d4	7.756	132	64541	24.452	ng/u1	0.00
15) 4-Methylphenol-d8	8.925	113	67678	22.572	ng/u1	0.00
21) Nitrobenzene-d5	9.401	128	34074	25.162	ng/u1	0.00
24) 2-Nitrophenol-d4	10.130	143	36081	23.962	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.670	165	56841	22.410	ng/u1	0.00
31) 4-Chloroaniline-d4	11.187	131	91757	23.888	ng/u1	0.00
46) Dimethylphthalate-d6	14.248	166	221173	27.254	ng/u1	0.00
49) Acenaphthylene-d8	14.548	160	266604	26.369	ng/u1	0.00
54) 4-Nitrophenol-d4	15.036	143	29892	20.315	ng/u1	0.00
60) Fluorene-d10	15.841	176	184875	25.716	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.958	200	27581	19.318	ng/u1	0.00
73) Anthracene-d10	17.697	188	311431	27.971	ng/u1	0.00
81) Pyrene-d10	19.971	212	355612	26.419	ng/u1	0.00
92) Benzo(a)pyrene-d12	25.059	264	292750	26.174	ng/u1	-0.01

Target Compounds Qvalue

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

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