

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG111521\
 Data File : BG051037.D
 Acq On : 15 Nov 2021 11:10
 Operator : CG/JU
 Sample : PB140632BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :

Quant Time: Nov 15 12:10:36 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG110321.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Nov 15 12:03:08 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.225	152	30983	20.000	ng/u1	0.00
20) Naphthalene-d8	11.051	136	133757	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.853	164	91303	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.596	188	214712	20.000	ng/u1	0.00
79) Chrysene-d12	21.897	240	206863	20.000	ng/u1	0.00
88) Perylene-d12	25.305	264	199565	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.578	96	5275	5.495	ng/uL	0.00
4) Pyridine-d5	4.006	84	76245	26.550	ng/u1	0.00
7) Phenol-d5	7.367	99	83799	25.353	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.538	67	53804	25.199	ng/u1	0.00
11) 2-Chlorophenol-d4	7.749	132	58338	25.467	ng/u1	0.00
15) 4-Methylphenol-d8	8.924	113	65373	25.123	ng/u1	0.00
21) Nitrobenzene-d5	9.400	128	31189	27.439	ng/u1	0.00
24) 2-Nitrophenol-d4	10.123	143	35539	28.119	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.669	165	56730	26.646	ng/u1	0.00
31) 4-Chloroaniline-d4	11.186	131	88486	27.445	ng/u1	0.00
46) Dimethylphthalate-d6	14.242	166	213670	30.589	ng/u1	0.00
49) Acenaphthylene-d8	14.547	160	250164	28.745	ng/u1	0.00
54) 4-Nitrophenol-d4	15.041	143	35853	28.308	ng/u1	0.00
60) Fluorene-d10	15.840	176	175125	28.301	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.957	200	30382	23.339	ng/u1	0.00
73) Anthracene-d10	17.696	188	306974	30.240	ng/u1	0.00
81) Pyrene-d10	19.976	212	382799	28.651	ng/u1	0.00
92) Benzo(a)pyrene-d12	25.064	264	318596	28.879	ng/u1	0.00

Target Compounds Qvalue

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

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