

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG121721\
 Data File : BG051619.D
 Acq On : 18 Dec 2021 7:12
 Operator : CG/JU
 Sample : PB141449BL
 Misc :
 ALS Vial : 34 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SBLK449

Quant Time: Dec 18 19:46:04 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG121421.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Dec 14 14:19:57 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.273	152	38060	20.000	ng/ul	-0.01
20) Naphthalene-d8	11.116	136	153664	20.000	ng/ul	# 0.00
38) Acenaphthene-d10	14.911	164	108003	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.655	188	254376	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.972	240	246523	20.000	ng/ul	#-0.01
88) Perylene-d12	25.468	264	237026	20.000	ng/ul	-0.03
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.603	96	5637	6.151	ng/uL	0.00
4) Pyridine-d5	4.032	84	21751	7.855	ng/ul	-0.01
7) Phenol-d5	7.404	99	107804	31.723	ng/ul	-0.01
9) Bis-(2-Chloroethyl)eth...	7.580	67	64005	31.683	ng/ul	-0.01
11) 2-Chlorophenol-d4	7.797	132	77811	32.922	ng/ul	-0.01
15) 4-Methylphenol-d8	8.966	113	84637	32.221	ng/ul	-0.01
21) Nitrobenzene-d5	9.442	128	40581	36.021	ng/ul	-0.01
24) 2-Nitrophenol-d4	10.177	143	43113	35.007	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.717	165	86411	32.355	ng/ul	-0.01
31) 4-Chloroaniline-d4	0.000	131	0	0.000	ng/ul	
46) Dimethylphthalate-d6	14.300	166	287935	33.308	ng/ul	-0.01
49) Acenaphthylene-d8	14.606	160	349284	32.534	ng/ul	0.00
54) 4-Nitrophenol-d4	15.064	143	45268	32.202	ng/ul	-0.01
60) Fluorene-d10	15.898	176	246243	31.794	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.998	200	42728	32.240	ng/ul	-0.01
73) Anthracene-d10	17.755	188	408188	33.639	ng/ul	0.00
81) Pyrene-d10	20.028	212	485528	32.724	ng/ul	-0.01
92) Benzo(a)pyrene-d12	25.227	264	435237	34.899	ng/ul	-0.03

Target Compounds Qvalue

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

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