

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120221\
 Data File : BG051313.D
 Acq On : 2 Dec 2021 21:50
 Operator : CG/JU
 Sample : PB141106BL
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SBLK106

Quant Time: Dec 03 00:02:35 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG112321.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Nov 24 06:04:50 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.196	152	32995	20.000	ng/u1	0.00
20) Naphthalene-d8	11.022	136	153893	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.823	164	104646	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.579	188	234800	20.000	ng/u1	0.00
79) Chrysene-d12	21.880	240	198536	20.000	ng/u1	0.00
88) Perylene-d12	25.276	264	195352	20.000	ng/u1	0.00

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.531	96	4972	5.237	ng/uL	-0.01
4) Pyridine-d5	3.965	84	74864	26.870	ng/u1	-0.01
7) Phenol-d5	7.356	99	92691	28.424	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.508	67	61966	30.256	ng/u1	0.00
11) 2-Chlorophenol-d4	7.726	132	68599	29.213	ng/u1	0.00
15) 4-Methylphenol-d8	8.907	113	76460	29.055	ng/u1	0.00
21) Nitrobenzene-d5	9.371	128	37173	28.615	ng/u1	0.00
24) 2-Nitrophenol-d4	10.099	143	43082	29.399	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.652	165	67463	27.133	ng/u1	0.00
31) 4-Chloroaniline-d4	11.163	131	104887	28.831	ng/u1	0.00
46) Dimethylphthalate-d6	14.218	166	240494	29.868	ng/u1	0.00
49) Acenaphthylene-d8	14.524	160	303226	29.865	ng/u1	0.00
54) 4-Nitrophenol-d4	15.052	143	31008	23.791	ng/u1	0.00
60) Fluorene-d10	15.816	176	208495	28.755	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.951	200	24618	16.991	ng/u1	0.00
73) Anthracene-d10	17.673	188	339300	30.215	ng/u1	0.00
81) Pyrene-d10	19.959	212	372248	30.987	ng/u1	0.00
92) Benzo(a)pyrene-d12	25.041	264	308930	29.610	ng/u1	0.00

Target Compounds Qvalue

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

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