

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG103123\
 Data File : BG059560.D
 Acq On : 31 Oct 2023 13:23
 Operator : MA/JU
 Sample : PB156468BL
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SBLK468

Quant Time: Nov 01 04:00:42 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG102723.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Sat Oct 28 03:56:42 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.360	152	77627	20.000	ng/u1	0.00
20) Naphthalene-d8	11.204	136	409545	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.988	164	227826	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.738	188	489177	20.000	ng/u1	0.00
79) Chrysene-d12	22.074	240	377311	20.000	ng/u1	0.00
88) Perylene-d12	25.681	264	416121	20.000	ng/u1	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.654	96	11797	5.686	ng/uL	0.00
4) Pyridine-d5	4.089	84	170147	29.347	ng/u1	0.00
7) Phenol-d5	7.467	99	269231	33.751	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.673	67	122794	33.514	ng/u1	0.00
11) 2-Chlorophenol-d4	7.873	132	189303	29.251	ng/u1	0.00
15) 4-Methylphenol-d8	9.036	113	196052	30.978	ng/u1	0.00
21) Nitrobenzene-d5	9.553	128	92920	31.909	ng/u1	0.00
24) 2-Nitrophenol-d4	10.276	143	104189	34.852	ng/u1	-0.01
28) 2,4-Dichlorophenol-d3	10.805	165	202050	30.677	ng/u1	0.00
31) 4-Chloroaniline-d4	11.345	131	326149	30.081	ng/u1	0.00
46) Dimethylphthalate-d6	14.383	166	645716	31.284	ng/u1	0.00
49) Acenaphthylene-d8	14.688	160	709132	29.275	ng/u1	0.00
54) 4-Nitrophenol-d4	15.152	143	109160	30.840	ng/u1	0.00
60) Fluorene-d10	15.975	176	516927	29.305	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	16.075	200	71607	30.911	ng/u1	0.00
73) Anthracene-d10	17.837	188	794154	29.966	ng/u1	0.00
81) Pyrene-d10	20.105	212	834927	31.836	ng/u1	0.00
92) Benzo(a)pyrene-d12	25.434	264	759483	30.995	ng/u1	0.00

Target Compounds Qvalue

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

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