

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG010424\
 Data File : BG060263.D
 Acq On : 5 Jan 2024 3:47
 Operator : MA/JU
 Sample : P1003-02
 Misc :
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 72-11978

Quant Time: Jan 05 04:24:58 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG122923.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Dec 30 03:14:04 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.936	152	166585	20.000	ng	0.00
21) Naphthalene-d8	10.738	136	707050	20.000	ng	0.00
39) Acenaphthene-d10	14.575	164	576185	20.000	ng	0.00
64) Phenanthrene-d10	17.319	188	1610996	20.000	ng	0.00
76) Chrysene-d12	21.585	240	1566627	20.000	ng	0.00
86) Perylene-d12	24.757	264	1745761	20.000	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.509	112	1525051	138.177	ng	0.00
7) Phenol-d6	7.101	99	1995004	121.741	ng	0.00
23) Nitrobenzene-d5	9.099	82	1431347	93.507	ng	0.00
42) 2,4,6-Tribromophenol	16.062	330	1203681	156.062	ng	0.00
45) 2-Fluorobiphenyl	13.194	172	3494881	87.176	ng	0.00
79) Terphenyl-d14	19.939	244	5595769	68.669	ng	0.00
Target Compounds						
77) Benzidine	19.563	184	892703	20.589	ng	Qvalue 99

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

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