

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120622\
 Data File : BG055908.D
 Acq On : 6 Dec 2022 18:05
 Operator : CG/JU
 Sample : N5884-01MSD
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 EO-3-120522MSD

Quant Time: Dec 07 05:11:31 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG111622.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Dec 07 05:04:10 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	0.000	152	0	0.000	ng	-8.16
21) Naphthalene-d8	0.000	136	0	0.000	ng	-10.99
39) Acenaphthene-d10	0.000	164	0	0.000	ng	-14.80
64) Phenanthrene-d10	0.000	188	0	0.000	ng	-17.55
76) Chrysene-d12	21.818	240	92	20.000	ng	#-0.01
86) Perylene-d12	25.138	264	166	20.000	ng	#-0.04
System Monitoring Compounds						
5) 2-Fluorophenol	0.000	112	0	0.000	ng	
7) Phenol-d6	0.000	99	0	0.000	ng	
23) Nitrobenzene-d5	0.000	82	0	0.000	ng	
42) 2,4,6-Tribromophenol	0.000	330	0	0.000	ng	
45) 2-Fluorobiphenyl	0.000	172	0	0.000	ng	
79) Terphenyl-d14	20.155	244	268	58.265	ng	0.03
Target Compounds						
						Qvalue
77) Benzidine	19.756	184	301	141.423	ng	# 1
78) Pyrene	19.950	202	531	85.270	ng	# 1
80) Butylbenzylphthalate	20.796	149	485	198.946	ng	# 56
81) Benzo(a)anthracene	21.801	228	945	149.134	ng	# 2
82) 3,3'-Dichlorobenzidine	21.712	252	187	83.986	ng	# 64
83) Chrysene	21.871	228	839	139.083	ng	# 5
84) Bis(2-ethylhexyl)phtha...	21.671	149	907	258.881	ng	# 60
85) Di-n-octyl phthalate	22.905	149	934	155.778	ng	# 1
87) Indeno(1,2,3-cd)pyrene	29.004	276	171	12.574	ng	# 1
88) Benzo(b)fluoranthene	24.051	252	108	9.990	ng	# 1
89) Benzo(k)fluoranthene	24.168	252	925	85.837	ng	# 1
90) Benzo(a)pyrene	25.003	252	344	37.081	ng	# 1
91) Dibenzo(a,h)anthracene	29.027	278	213	19.167	ng	# 1
92) Benzo(g,h,i)perylene	30.214	276	138	12.321	ng	# 17

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

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