

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG112423\
 Data File : BG059874.D
 Acq On : 25 Nov 2023 1:47
 Operator : MA/JU
 Sample : 05408-07
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 712-PLAINFIELD-AVE

Quant Time: Nov 25 03:47:56 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG111623.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Nov 24 15:11:07 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.323	152	26117	20.000	ng	0.00
21) Naphthalene-d8	11.166	136	154822	20.000	ng	# 0.00
39) Acenaphthene-d10	14.950	164	118356	20.000	ng	0.00
64) Phenanthrene-d10	17.700	188	269250	20.000	ng	# 0.00
76) Chrysene-d12	22.018	240	208201	20.000	ng	#-0.02
86) Perylene-d12	25.585	264	230557	20.000	ng	-0.03
System Monitoring Compounds						
5) 2-Fluorophenol	5.814	112	133271	85.888	ng	0.00
7) Phenol-d6	7.441	99	211207	91.851	ng	0.00
23) Nitrobenzene-d5	9.515	82	232225	57.286	ng	0.00
42) 2,4,6-Tribromophenol	16.437	330	140684	104.762	ng	0.00
45) 2-Fluorobiphenyl	13.570	172	495244	56.323	ng	0.00
79) Terphenyl-d14	20.273	244	812347	57.599	ng	0.00
Target Compounds						
						Qvalue
75) Fluoranthene	19.733	202	61231	3.730	ng	99
78) Pyrene	20.097	202	53037	3.744	ng	98
81) Benzo(a)anthracene	22.001	228	30416	2.135	ng	# 87
83) Chrysene	22.071	228	31691	2.427	ng	# 85
88) Benzo(b)fluoranthene	24.433	252	40547	3.209	ng	# 98
90) Benzo(a)pyrene	25.409	252	33183	2.460	ng	# 87

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

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