

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG020723\
 Data File : BG056572.D
 Acq On : 7 Feb 2023 18:02
 Operator : CG/JU
 Sample : 01462-15
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 COMP-3

Quant Time: Feb 08 03:06:48 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG012723.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Feb 08 03:02:11 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.269	152	33674	20.000	ng	0.00
21) Naphthalene-d8	11.101	136	129871	20.000	ng	0.00
39) Acenaphthene-d10	14.885	164	99380	20.000	ng	0.00
64) Phenanthrene-d10	17.629	188	240947	20.000	ng	0.00
76) Chrysene-d12	21.918	240	217256	20.000	ng	0.00
86) Perylene-d12	25.332	264	235696	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.790	112	226779	123.887	ng	0.00
7) Phenol-d6	7.417	99	311185	114.734	ng	0.00
23) Nitrobenzene-d5	9.444	82	163703	71.578	ng	0.00
42) 2,4,6-Tribromophenol	16.371	330	146082	110.568	ng	0.00
45) 2-Fluorobiphenyl	13.516	172	526142	73.401	ng	0.00
79) Terphenyl-d14	20.202	244	902039	72.924	ng	0.00
Target Compounds						
75) Fluoranthene	19.662	202	50423	3.168	ng	Qvalue 100
78) Pyrene	20.026	202	48384	3.131	ng	98
88) Benzo(b)fluoranthene	24.239	252	33799	2.310	ng	# 93

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

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