

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG012424\
 Data File : BG060422.D
 Acq On : 24 Jan 2024 19:39
 Operator : MA/JU
 Sample : P1234-01
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 293

Quant Time: Jan 25 00:07:55 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG010824.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jan 09 04:07:12 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.930	152	154566	20.000	ng	0.00
21) Naphthalene-d8	10.726	136	671479	20.000	ng	0.00
39) Acenaphthene-d10	14.563	164	564932	20.000	ng	0.00
64) Phenanthrene-d10	17.313	188	1561862	20.000	ng	0.00
76) Chrysene-d12	21.572	240	1423767	20.000	ng	0.00
86) Perylene-d12	24.739	264	1604071	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.503	112	1173389	124.864	ng	0.00
7) Phenol-d6	7.089	99	1444987	103.842	ng	0.00
23) Nitrobenzene-d5	9.087	82	870814	61.238	ng	0.00
42) 2,4,6-Tribromophenol	16.055	330	974180	117.205	ng	0.00
45) 2-Fluorobiphenyl	13.182	172	2320740	57.114	ng	0.00
79) Terphenyl-d14	19.927	244	4004725	64.888	ng	0.00
Target Compounds						
						Qvalue
71) Phenanthrene	17.354	178	386298	5.117	ng	92
75) Fluoranthene	19.369	202	447812	4.939	ng	87
78) Pyrene	19.727	202	350374	3.854	ng	# 85

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

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