

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG123022\
 Data File : BG056178.D
 Acq On : 30 Dec 2022 19:58
 Operator : CG/JU
 Sample : N6237-01
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 TR-05-122922

Quant Time: Dec 31 00:20:30 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG122722.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Dec 28 05:20:14 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.340	152	55566	20.000	ng	0.00
21) Naphthalene-d8	11.166	136	203090	20.000	ng	0.00
39) Acenaphthene-d10	14.944	164	152717	20.000	ng	0.00
64) Phenanthrene-d10	17.676	188	359062	20.000	ng	0.00
76) Chrysene-d12	21.971	240	348727	20.000	ng	0.00
86) Perylene-d12	25.432	264	374458	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.861	112	216426	69.544	ng	0.00
7) Phenol-d6	7.471	99	294742	61.671	ng	0.00
23) Nitrobenzene-d5	9.504	82	151914	38.262	ng	0.00
42) 2,4,6-Tribromophenol	16.419	330	112810	58.326	ng	0.00
45) 2-Fluorobiphenyl	13.569	172	428734	38.767	ng	0.00
79) Terphenyl-d14	20.250	244	710564	36.494	ng	0.00
Target Compounds						
						Qvalue
71) Phenanthrene	17.723	178	51531	2.769	ng	96
75) Fluoranthene	19.709	202	127454	5.251	ng	100
78) Pyrene	20.068	202	119651	4.868	ng	98
81) Benzo(a)anthracene	21.954	228	62400	2.548	ng	99
83) Chrysene	22.024	228	63941	2.787	ng	94
88) Benzo(b)fluoranthene	24.322	252	87305	3.670	ng	100
90) Benzo(a)pyrene	25.262	252	60678	2.615	ng	# 98

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

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