

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG011323\
 Data File : BG056282.D
 Acq On : 13 Jan 2023 17:57
 Operator : CG/JU
 Sample : 01135-01
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 WC-2

Quant Time: Jan 16 00:48:24 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG122722.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jan 12 03:38:40 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.319	152	38024	20.000	ng	0.00
21) Naphthalene-d8	11.151	136	147758	20.000	ng	# 0.00
39) Acenaphthene-d10	14.929	164	125180	20.000	ng	0.00
64) Phenanthrene-d10	17.667	188	317813	20.000	ng	0.00
76) Chrysene-d12	21.956	240	315126	20.000	ng	# 0.00
86) Perylene-d12	25.399	264	342157	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.834	112	278526	130.788	ng	0.00
7) Phenol-d6	7.449	99	401966	122.908	ng	0.00
23) Nitrobenzene-d5	9.488	82	212656	73.618	ng	0.00
42) 2,4,6-Tribromophenol	16.409	330	184338	116.273	ng	0.00
45) 2-Fluorobiphenyl	13.554	172	630117	69.510	ng	0.00
79) Terphenyl-d14	20.240	244	1255877	71.379	ng	0.00
Target Compounds						
						Qvalue
71) Phenanthrene	17.708	178	52412	3.182	ng	97
75) Fluoranthene	19.694	202	93464	4.350	ng	97
78) Pyrene	20.058	202	90700	4.083	ng	100
81) Benzo(a)anthracene	21.932	228	50070	2.262	ng	99
83) Chrysene	22.009	228	42997	2.074	ng	96
88) Benzo(b)fluoranthene	24.294	252	52369	2.409	ng	# 88

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

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