

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG123022\
 Data File : BG056179.D
 Acq On : 30 Dec 2022 20:39
 Operator : CG/JU
 Sample : N6235-01 5X
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 OR-2-122922

Quant Time: Dec 31 00:20:46 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	8.336	152	64555	20.000	ng	0.00
21) Naphthalene-d8	11.168	136	228780	20.000	ng	0.00
39) Acenaphthene-d10	14.946	164	170666	20.000	ng	0.00
64) Phenanthrene-d10	17.678	188	395700	20.000	ng	0.00
76) Chrysene-d12	21.979	240	372519	20.000	ng	0.00
86) Perylene-d12	25.439	264	405064	20.000	ng	0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.856	112	66239	18.321	ng	0.00
7) Phenol-d6	7.466	99	94668	17.050	ng	0.00
23) Nitrobenzene-d5	9.505	82	47693	10.663	ng	0.00
42) 2,4,6-Tribromophenol	16.420	330	34054	15.755	ng	0.00
45) 2-Fluorobiphenyl	13.571	172	142660	11.543	ng	0.00
79) Terphenyl-d14	20.251	244	228817	11.001	ng	0.00
Target Compounds						
						Qvalue
71) Phenanthrene	17.725	178	94382	4.603	ng	99
75) Fluoranthene	19.711	202	171376	6.407	ng	98
78) Pyrene	20.069	202	216568	8.248	ng	99
81) Benzo(a)anthracene	21.955	228	98802	3.776	ng	96
83) Chrysene	22.026	228	99257	4.050	ng	97
88) Benzo(b)fluoranthene	24.329	252	117013	4.547	ng	95
90) Benzo(a)pyrene	25.269	252	96879	3.859	ng	99
92) Benzo(g,h,i)perylene	30.663	276	59848	2.499	ng	# 92

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

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