

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG080222\
 Data File : BG054482.D
 Acq On : 2 Aug 2022 14:03
 Operator : CG/JU
 Sample : N3894-03
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 RVE-22

Quant Time: Aug 03 04:13:35 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG071522.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jul 28 04:30:50 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.984	152	16946	20.000	ng	# 0.00
21) Naphthalene-d8	10.792	136	65645	20.000	ng	# 0.00
39) Acenaphthene-d10	14.623	164	51672	20.000	ng	0.00
64) Phenanthrene-d10	17.367	188	135053	20.000	ng	# 0.00
76) Chrysene-d12	21.632	240	151813	20.000	ng	# 0.00
86) Perylene-d12	24.840	264	161077	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.569	112	91247	112.424	ng	0.00
7) Phenol-d6	7.179	99	124432	111.885	ng	0.00
23) Nitrobenzene-d5	9.147	82	83294	69.098	ng	0.00
42) 2,4,6-Tribromophenol	16.115	330	122764	113.155	ng	0.00
45) 2-Fluorobiphenyl	13.242	172	261493	71.390	ng	0.00
79) Terphenyl-d14	19.976	244	573914	74.996	ng	0.00
Target Compounds						
						Qvalue
75) Fluoranthene	19.423	202	35184	3.838	ng	95
78) Pyrene	19.782	202	46835	5.377	ng	99
83) Chrysene	21.679	228	32375	3.556	ng	95
88) Benzo(b)fluoranthene	23.824	252	39700	4.168	ng	93
90) Benzo(a)pyrene	24.694	252	27056	3.326	ng	99

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

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