

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG080122\
 Data File : BG054469.D
 Acq On : 1 Aug 2022 18:02
 Operator : CG/JU
 Sample : N3896-05
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 RVE-62

Manual Integrations
 APPROVED

Reviewed By : Christian Giraldo 08/02/2022
 Supervised By : Jagrut Upadhyay 08/02/2022

Quant Time: Aug 02 05:46:19 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	7.978	152	16355	20.000	ng	0.00
21) Naphthalene-d8	10.793	136	66224	20.000	ng	# 0.00
39) Acenaphthene-d10	14.618	164	55246	20.000	ng	0.00
64) Phenanthrene-d10	17.367	188	138163	20.000	ng	# 0.00
76) Chrysene-d12	21.633	240	151241	20.000	ng	# 0.00
86) Perylene-d12	24.835	264	162000	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.564	112	101094	129.057	ng	0.00
7) Phenol-d6	7.179	99	139965	130.400	ng	0.00
23) Nitrobenzene-d5	9.148	82	93054	76.520	ng	0.00
42) 2,4,6-Tribromophenol	16.116	330	142987	123.269	ng	0.00
45) 2-Fluorobiphenyl	13.243	172	306404	78.240	ng	0.00
79) Terphenyl-d14	19.976	244	640355	83.994	ng	0.00
Target Compounds						Qvalue
75) Fluoranthene	19.418	202	25131	2.679	ng	96
78) Pyrene	19.782	202	26287	3.029	ng	94
88) Benzo(b)fluoranthene	23.825	252	41213m	4.302	ng	
90) Benzo(a)pyrene	24.688	252	24425	2.985	ng	96

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

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