

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG102922\
 Data File : BG055374.D
 Acq On : 29 Oct 2022 20:53
 Operator : CG/JU
 Sample : N5313-01
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 COMPOSITE-SW-QUARTER

Quant Time: Oct 31 02:45:26 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG102422.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Oct 29 04:52:38 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.261	152	34879	20.000	ng	0.00
21) Naphthalene-d8	11.093	136	148551	20.000	ng	# 0.00
39) Acenaphthene-d10	14.871	164	97916	20.000	ng	0.00
64) Phenanthrene-d10	17.609	188	210401	20.000	ng	# 0.00
76) Chrysene-d12	21.881	240	199568	20.000	ng	# 0.00
86) Perylene-d12	25.277	264	217525	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.788	112	246832	123.913	ng	0.00
7) Phenol-d6	7.410	99	323617	112.137	ng	0.00
23) Nitrobenzene-d5	9.437	82	197442	67.869	ng	0.00
42) 2,4,6-Tribromophenol	16.352	330	143561	134.871	ng	0.00
45) 2-Fluorobiphenyl	13.502	172	460870	77.039	ng	0.00
79) Terphenyl-d14	20.183	244	683848	71.195	ng	0.00
Target Compounds						
						Qvalue
71) Phenanthrene	17.650	178	28092	2.504	ng	99
75) Fluoranthene	19.636	202	64116	4.904	ng	97
78) Pyrene	20.001	202	61436	4.491	ng	98
81) Benzo(a)anthracene	21.863	228	34464	2.572	ng	98
83) Chrysene	21.928	228	30307	2.381	ng	94
88) Benzo(b)fluoranthene	24.190	252	41451	3.168	ng	97
90) Benzo(a)pyrene	25.106	252	31406	2.799	ng	98

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

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