

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG031022\
 Data File : BG052635.D
 Acq On : 10 Mar 2022 15:41
 Operator : CG/JU
 Sample : N1845-10 2X
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SAMPLE-10

Manual Integrations
 APPROVED

Reviewed By : Christian Giraldo 03/11/2022
 Supervised By : Jagrut Upadhyay 03/11/2022

Quant Time: Mar 11 03:45:00 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG030922.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Mar 10 06:37:54 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.175	152	45221	20.000	ng	0.00
21) Naphthalene-d8	11.012	136	169803	20.000	ng	0.00
39) Acenaphthene-d10	14.825	164	108111	20.000	ng	0.00
64) Phenanthrene-d10	17.580	188	220702	20.000	ng	0.00
76) Chrysene-d12	21.886	240	190349	20.000	ng	0.00
86) Perylene-d12	25.316	264	186913	20.000	ng	# 0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.707	112	91316	34.885	ng	0.00
7) Phenol-d6	7.329	99	148912	39.306	ng	0.00
23) Nitrobenzene-d5	9.350	82	102020	27.236	ng	0.00
42) 2,4,6-Tribromophenol	16.317	330	42795	27.910	ng	0.00
45) 2-Fluorobiphenyl	13.450	172	229907	28.704	ng	0.00
79) Terphenyl-d14	20.170	244	312906	29.089	ng	0.00
Target Compounds						
						Qvalue
71) Phenanthrene	17.621	178	23936	2.067	ng	98
75) Fluoranthene	19.624	202	44296	3.064	ng	100
78) Pyrene	19.988	202	41854m	3.290	ng	
83) Chrysene	21.933	228	25179	2.112	ng	# 89
88) Benzo(b)fluoranthene	24.224	252	27108	2.349	ng	# 69

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

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