

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG110119\
 Data File : BG043478.D
 Acq On : 1 Nov 2019 23:06
 Operator : JU
 Sample : K5541-17 5X
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 18-48-(0-2.5)

Manual Integrations
 APPROVED

mohammad
 11/4/2019 11:50:02 AM

Quant Time: Nov 02 00:32:08 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG102119.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Oct 21 16:37:43 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.02	152	53462	20.00	ng	0.00
21) Naphthalene-d8	10.82	136	196693	20.00	ng	0.00
39) Acenaphthene-d10	14.64	164	136876	20.00	ng	0.00
64) Phenanthrene-d10	17.39	188	292571	20.00	ng	-0.02
76) Chrysene-d12	21.65	240	301241	20.00	ng	-0.02
87) Perylene-d12	24.86	264	365337	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.61	112	42207	14.47	ng	0.00
7) Phenol-d6	7.20	99	58471	13.93	ng	0.00
23) Nitrobenzene-d5	9.18	82	35279	9.25	ng	0.00
42) 2,4,6-Tribromophenol	16.14	330	31864	12.23	ng	0.00
45) 2-Fluorobiphenyl	13.27	172	90103	9.10	ng	0.00
79) Terphenyl-d14	20.00	244	146608	9.13	ng	-0.01
Target Compounds						
75) Fluoranthene	19.44	202	70933	3.632	ng	99
78) Pyrene	19.80	202	68803	3.690	ng	98
81) Benzo(a)anthracene	21.64	228	45690	2.421	ng	97
83) Chrysene	21.70	228	40499m	2.160	ng	
88) Benzo(b)fluoranthene	23.85	252	60173	2.908	ng	# 82
90) Benzo(a)pyrene	24.70	252	45146	2.285	ng	# 88

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

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