

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF102918\
 Data File : BF110276.D
 Acq On : 29 Oct 2018 21:45
 Operator : JU/SJ
 Sample : J5665-03
 Misc :
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 DL-07A

Manual Integrations
 APPROVED

Sohil
 10/30/2018 4:42:45 PM

Quant Time: Oct 30 12:28:38 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF102318.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Oct 29 17:29:00 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.97	152	117870	20.00	ng	0.00
21) Naphthalene-d8	8.25	136	485420	20.00	ng	0.00
39) Acenaphthene-d10	10.01	164	221581	20.00	ng	0.00
64) Phenanthrene-d10	11.50	188	360774	20.00	ng	0.00
76) Chrysene-d12	14.15	240	235774	20.00	ng	0.00
87) Perylene-d12	15.69	264	200369	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.60	112	693831	95.86	ng	0.01
7) Phenol-d6	6.59	99	916707	99.02	ng	0.00
23) Nitrobenzene-d5	7.53	82	576347	70.42	ng	0.00
42) 2,4,6-Tribromophenol	10.80	330	161812	73.72	ng	0.00
45) 2-Fluorobiphenyl	9.33	172	965320	68.41	ng	0.00
79) Terphenyl-d14	13.09	244	640111	56.46	ng	0.00
Target Compounds						
10) Phenol	6.60	94	74341	7.512	ng	# 64
22) Acetophenone	7.35	105	25175	2.251	ng	# 61
31) Naphthalene	8.27	128	247741	11.073	ng	99
37) 2-Methylnaphthalene	8.96	142	248957	16.819	ng	98
38) 1-Methylnaphthalene	9.06	142	199149m	13.922	ng	
46) 1,1'-Biphenyl	9.43	154	41009	2.047	ng	98
50) Dimethylphthalate	9.72	163	202517	12.381	ng	99
55) Dibenzofuran	10.22	168	65040	3.308	ng	# 96
71) Phenanthrene	11.53	178	168553	8.929	ng	100
78) Pyrene	12.96	202	35928m	2.126	ng	
83) Chrysene	14.18	228	41909	3.156	ng	# 87
88) Benzo(b)fluoranthene	15.23	252	24346m	2.104	ng	

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

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