

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF102618\
 Data File : BF110227.D
 Acq On : 26 Oct 2018 23:13
 Operator : JU/SJ
 Sample : J5675-01
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TP-5

Manual Integrations
 APPROVED

Sohil
 10/29/2018 11:56:49 AM

Quant Time: Oct 27 00:55:37 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF102318.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Oct 23 15:06:18 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.97	152	126109	20.00	ng	-0.01
21) Naphthalene-d8	8.25	136	478692	20.00	ng	-0.02
39) Acenaphthene-d10	10.01	164	186467	20.00	ng	-0.02
64) Phenanthrene-d10	11.50	188	286482	20.00	ng	-0.01
76) Chrysene-d12	14.16	240	197700	20.00	ng	0.00
87) Perylene-d12	15.69	264	133458	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.59	112	727481	93.94	ng	0.00
7) Phenol-d6	6.59	99	904816	91.35	ng	-0.01
23) Nitrobenzene-d5	7.53	82	531312	65.83	ng	-0.02
42) 2,4,6-Tribromophenol	10.80	330	143635	77.76	ng	-0.01
45) 2-Fluorobiphenyl	9.33	172	832043	70.07	ng	-0.01
79) Terphenyl-d14	13.09	244	585129	61.55	ng	-0.01
Target Compounds						
10) Phenol	6.60	94	65691	6.204	ng	# 64
50) Dimethylphthalate	9.72	163	154890	11.253	ng	99
75) Fluoranthene	12.72	202	54913	3.610	ng	97
78) Pyrene	12.95	202	54549	3.849	ng	98
81) Benzo(a)anthracene	14.14	228	36513	3.114	ng	# 83
83) Chrysene	14.18	228	33109	2.973	ng	96
86) Indeno(1,2,3-cd)pyrene	17.26	276	29274	3.169	ng	98
88) Benzo(b)fluoranthene	15.23	252	45979m	5.966	ng	
90) Benzo(a)pyrene	15.62	252	32183	4.538	ng	# 79
92) Benzo(g,h,i)perylene	17.74	276	28675	4.756	ng	# 83

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

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